

Memorandum

To: Stephanie Vaughn, EPA Region 2

Elizabeth Buckrucker, USACE

From: Frank Tsang and Scott Kirchner (CDM Smith)

Date: January 24, 2014

Subject: 2012 Background Sediment Split Sample Data Comparison for the Lower

Passaic River Study Area

At the request of the United State Environmental Protection Agency (EPA) and the United States Army Corps of Engineers (USACE), CDM Federal Programs Corporation (CDM Smith) collected oversight split background sediment samples as part of the Lower Passaic River (LPR) Restoration Project remedial investigation conducted by the Cooperating Parties Group (CPG). This memorandum presents the comparison of the EPA oversight team's split sample results to the CPG's sample results and discusses the differences in the datasets. In this document, samples will be referred to as either CPG samples or EPA samples for clarity.

The evaluation was conducted in two ways: (1) on an individual chemical basis and (2) on a chemical group basis. Normally, only the first of these (individual chemical basis) would be conducted, consistent with previous comparisons. However, due to the small number of data pairs available for the individual chemicals, which resulted in insufficient statistical power to discern differences, it was determined that the second method (chemical group basis) should also be conducted. The split sample data comparison for the 2012 Lower Passaic background sediment sampling oversight is based primarily on the second method except for total organic carbon (TOC) and mercury.

- <u>Dioxins/Furans</u>: For dioxins and furans as a group, there were statistically significant differences between the results of the EPA and CPG background sediment samples. Overall, the CPG results were higher than the EPA results.
- <u>Pesticides</u>: The analytical results for pesticides as a group were comparable for EPA and CPG split samples. Where differences existed, CPG sample results tended to be lower than EPA samples results.
- <u>Polychlorinated Biphenyls (PCBs)</u>: The analytical results for PCBs as a group were comparable for EPA and CPG split samples. Where differences existed, CPG sample results tended to be higher than EPA samples results.

- Polycyclic Aromatic Hydrocarbons (PAHs): The analytical results for PAHs as a group were comparable for EPA and CPG split samples. Where differences existed, CPG sample results tended to be higher than EPA samples results.
- <u>TOC</u>: Though rigorous statistical testing was not possible due to the small sample size, the CPG results tended to be higher than the EPA results.
- Metals (excluding mercury): The analytical results for metals as a group were comparable for EPA and CPG split samples. Where differences existed, CPG sample results tended to be higher than EPA sample results.
- Mercury: The mercury results were comparable for CPG and EPA results.

Oversight Program Summary

Oversight was conducted in accordance with the Final Quality Assurance Project Plan (QAPP), Addendum No. 12, Collection of Background Surface Sediment Samples. The split sample program consisted of four background sediment split samples collected from the study area, therefore results from four split sample pairs were compared for each of the analytical methods evaluated.

Data Comparison Methodology

The CPG and EPA split sample data were evaluated for potential difference by plotting selected analytes (Table 1) for the following parameters. For each of the following three plots, data are plotted and evaluated only for the cases where both sample pairs are detected:

- <u>Line Plot of Absolute Concentrations</u>: The absolute concentrations measured by both analytical programs for the detected paired samples were plotted against the same axes. These graphs depict the relative magnitudes and patterns of concentrations.
- Bivariate Scatter Plot: CPG sample concentration was plotted as a function of EPA sample concentration for each detected pair. The bivariate plot illustrates the relationship between EPA and CPG data. Also included on each graph is a line which depicts a 1:1 ratio of EPA to CPG concentration. The bivariate plot can be used to identify potential systematic bias when data points fall consistently above or below the 1:1 line.
- <u>Percent Difference Plot</u>: The percent difference (%D) was defined as the difference between EPA and CPG concentration for detected data pairs, divided by the EPA sample concentration according to Equation 1.

$$\% D = \frac{(R_{USEPA} - R_{CPG})}{(R_{USEPA})} (100)$$
 (Equation 1)

Consequently, a negative %D indicates a CPG result that is higher than the EPA result, while a positive %D indicates a CPG result that is lower than the EPA result. This plot provides a visual

indication of the extent of positive and negative differences between the two datasets. The red dashed lines on the plot correspond to 40%D and -67%D. These criteria correspond to 50% relative percent difference (RPD) (the CPG's field duplicate acceptance criterion), converted to %D values. %D is commonly used when one of the two values is known or accepted, whereas RPD is more commonly used when both values are uncertain. The sample data in these graphs was represented with the EPA result as the known value and the CPG result as the unknown value.

In addition to the preparation of data comparison plots, the tests described below were also conducted for CPG and EPA data pairs where both results were detected.

- Average Ratio: The ratios of the CPG results to EPA results were calculated for each sample pair. The average ratio and standard deviation were calculated. An average ratio above one indicates that the CPG results were detected higher than the EPA results, while an average ratio below one indicates that the CPG results were detected lower than the EPA results.
- Percent Difference: The calculated %D values were evaluated against the acceptance criteria of greater than or equal to -67% or less than or equal to 40% (equivalent to less than or equal to 50% RPD). Since RPD is evaluated for results at least five times the sample reporting limit, only sample data pairs where both results were at least five times the sample reporting limit were included in the %D evaluation.
- Wilcoxon Signed Rank Test: The Wilcoxon Signed Rank test was used to calculate *p*-values for all detected sample pairs. The *p*-value is an indicator of the presence of a difference between the datasets. A *p*-value of less than 0.05 indicates a statistically significant difference between the two data sets. In addition to the regular Wilcoxon test for the detected data pairs, the Paired Prentice-Wilcoxon (PPW) test, a modified version of the Wilcoxon test, was utilized to allow inclusion of non-detected values. Further details on the PPW test are provided later in this memorandum.

The test criteria listed above were applied to two types of data groupings: (1) to the individual chemicals and (2) to the chemical groups as a whole. The data comparison plots associated with these two types of groupings are provided in Figures 1 through 48 for the individual chemicals, and in Figures 49 through 53 for the chemical groups. Results for the three statistical tests are presented in Table 1 for the individual chemicals and in Table 2 for the chemical groups. Other relevant information about the datasets is also provided in Tables 1 and 2. The numbers of split sample pairs are listed for each compound/group along with the number of pairs which had detected results for both samples. The average ratio of CPG sample results to EPA sample results are reported along with the standard deviation of the ratios. The %D results are summarized by reporting the percentage of data pairs that exceeded the acceptance criteria. Also included are the *p*-values calculated by the Wilcoxon Signed Rank test and PPW test.

An overall evaluation of the split sample data is based on the result of the three statistical tests, where each compound has a rating of "Same" or "Different". The datasets are considered comparable or "Same" if the criteria for at least two of the three statistical tests were met. When only the average ratio and Wilcoxon Signed Rank test (and/or PPW test) results were available, the overall ranking for each compound was considered comparable or "Same" if criteria for both tests were met. The criteria for each test are:

- Average Ratio: Average ratio of CPG to EPA within 0.70 to 1.30.
- Percent Difference: %D within 40 to -67% for the majority of the sample pairs, where no more than 16% of the data pairs exceed the acceptance criteria.
- <u>Wilcoxon Signed Rank Test (and/or PPW test)</u>: *p*-values greater than or equal to 0.05 are within acceptance limits, indicating there is no significant difference between the data sets.

In summary, the split sample comparison of the 2012 LPR background sediment consisted of four sample pairs. The split comparison was therefore limited to four sample pairs analyzed for dioxins and furans, organochlorine pesticides, PCBs, PAHs, TOC, metals, and mercury.

As discussed previously, the results presented in Table 1 (individual chemical basis) are limited due to the small sample size of the data pairs for the individual organic compounds and metals available for the evaluation. Specifically, with only a maximum of four data pairs, the statistical power of the Wilcoxon Signed Rank test (and/or PPW test) is insufficient to discern a real difference for the paired data. While the nature of the actual reduction in statistical power resulting from testing with four or less four data pairs was not investigated, such reduction is inherent in any statistical-based comparison test and a good general rule-of-thumb is a minimum of 8 to 10 samples or data pairs for robust statistical testing.

Due to the limitation for individual chemical data sets, a second evaluation (Table 2) was conducted whereby the data were tested as chemical groups: dioxins/furans, pesticides, PCBs, PAHs, and metals (excluding mercury). This testing was conducted to increase the number of sample pairs and thus increase statistical power. To account for differences in concentration ranges between the various compounds/metals within a group, each data pair was first scaled between 0 and 1 according to Equation 2,

$$y_i = \frac{R_i}{maximum(R_{USEPA,R_{CPG}})}$$
 (Equation 2)

where

y_i = scaled data value for each individual data R_i = data value in the original concentration units

maximum (R_{EPA} , R_{CPG}) = the maximum concentration in the split sample pair

The paired Wilcoxon test was then conducted on the scaled data pairs. It is noted that conducting the tests on chemical groups rather than on individual compounds/metals results in the loss of the ability to evaluate differences for the individual compounds/metals. In such a case where the test discerns a statistical difference for a chemical group, it is not possible to determine which individual compounds/metals within the group may be responsible for the difference.

The standard paired Wilcoxon test was conducted only on data pairs where both results were detected values. This protocol is consistent with previous split sample comparisons conducted at the LPR study area. However, it has been noted that the elimination of data pairs containing nondetected values is essentially equivalent to ignoring potentially substantial information contained within these nondetect-containing data pairs, and may lead to biased results, and that therefore an attempt should be made to include them in the analyses. Therefore, in addition to the standard paired Wilcoxon test conducted on the detected data pairs only, a modified version of the test called the PPW test was conducted that allows inclusion of the left-censored (non-detected) data pairs. The PPW test relies on survival analysis computations as detailed in O'Brien and Fleming (1987) and is considered the standard test for the case of censored matched pairs (Helsel 2005).

Below are some notable observations from Tables 1 and 2 and Figures 1 through 53. As mentioned earlier, the statistical power of the Wilcoxon Signed Rank test and/or the PPW test is insufficient to discern a real difference for the paired data for individual compounds due to the small sample size (<5). Thus, only results for the average ratio and %D are discussed below for individual compounds.

Dioxins/Furans

There was a high degree of variance between the CPG and EPA results for dioxin congener analysis. Of the six dioxin/furan congeners and total tetrachlorodibenzo-p-dioxin (TCDD) evaluated, only 2,3,7,8-TCDD and 2,3,7,8-tetrachlorodibenzofuran (2,3,7,8-TCDF)met the criteria for average ratios and %D, indicating that the split samples were comparable for these two congeners. For all other dioxin/furan congeners and total TCDD, the average ratio and %D exceeded acceptance criteria (Table 1). Differences are observed between the two data sets for 1,2,3,4,6,7,8-HpCDD, 1,2,3,4,6,7,8-HpCDF, OCDD, OCDF, and total TCDD based on the average ratio and %D. The CPG results were distinctly higher than the EPA results.

The differences between the dioxin data sets were further confirmed by the group testing. There was a statistically significant difference between the EPA and CPG samples. The average ratio and %D also exceeded acceptance criteria (Table 2). As shown in Figures 1 through 7 and Figure 49, overall the CPG results were high in comparison to the EPA results.

Pesticides

The comparison of analytical results between the EPA and the CPG showed comparable results for the majority of the pesticides. The average ratios criterion was met for all pesticides except 2,4'-dichlorodiphenyldichloroethane (DDD). The %D criterion was met for all pesticides except 2,4'-DDD and 4,4'-DDD. All pesticides data were comparable except 2,4'-DDD, where the CPG results were lower than the EPA results. From examination of individual sample pairs, the most marked discrepancy was observed for the 2,4'-DDD results for sample UPRT18C, where the CPG result was nearly 50 times lower than the EPA result.

The analytical results for pesticides as a group were comparable for EPA and CPG split samples. The PPW test did indicate that the data sets were different, but all other data evaluation indicated that the data sets were comparable.

Polychlorinated Biphenyls

The comparison of analytical results between the EPA and the CPG showed comparable results for the majority of the individual PCB congeners. Eight of the 11 PCB congeners evaluated and total PCB results were within the acceptable average ratios criterion. However, seven congeners and total PCB had %D exceeding the acceptance criterion. Differences were found between the data sets for 3,3',4,4'-tetrachlorobiphenyl (PCB 77), 3,4,4',5-tetrachlorobiphenyl (PCB 81), and 2,3,4,4',5-pentachlorobiphenyl (PCB 114) based on the average ratio and %D. The CPG results were higher than the EPA results for these congeners.

For PCBs as a group, the analytical data between the EPA and CPG samples were comparable, although the %D exceeded the acceptance criterion.

Polycyclic Aromatic Hydrocarbons

The split sample data were comparable for all individual PAH compounds except fluoranthene. All of the PAHs had average ratios within criterion except fluoranthene. Five of nine PAHs had %D within the criterion. The CPG results for fluoranthene were higher than the EPA results.

All three criteria indicated that there was no statistically significant difference between the EPA and CPG samples for PAHs as a group.

Total Organic Carbon

Differences were found between the two data sets for the TOC results, based on the average ratio and %D. The CPG results were higher than the EPA results (Table 1 and Figure 37). However, due to small sample size, the Wilcoxon Signed Rank test and the PPW test are insufficient to discern a real difference for the paired data.

Metals (excluding mercury)

The comparison of analytical results between the EPA and the CPG showed comparable results for the majority of metals. The average ratios were within the criterion for all metals except cadmium and lead. Six of ten metals had %D exceeding the criterion. Overall, the metal data are comparable except cadmium and lead. The CPG results were higher than the EPA results for these two metals. A notable discrepancy was noted for cadmium in sample UPRT20A. The CPG cadmium result was nearly 30 times higher than EPA result.

For metals as a group, there was no statistically significant difference between the EPA samples and CPG samples, though the %D exceeded the criterion.

Mercury

The mercury results were comparable for CPG and EPA results, based on the average ratio and %D.

Recommendations

Due to the small sample size (<5), the statistical power of the Wilcoxon signed rank test is greatly limited to identify a difference between the CPG and EPA data sets. In order to obtain meaningful statistical results for individual chemicals, at least 8 to 10 (preferably 15 to 20) split samples are recommended to collect during each oversight task, even if the number of split samples exceeds 10% of the number of the total samples.

Ignoring nondetects in the statistical comparison testing is an unnecessary practice, and hence should be avoided in future split sample comparisons. In this particular case, the PPW test should be conducted in lieu of the standard Wilcoxon test on detected values only.

Attachments

Table 1: Lower Passaic River 2012 Background Sediment Sampling Comparison Summary

by Individual Compounds

Table 2: Lower Passaic River 2012 Background Sediment Sampling Comparison Summary

by Chemical Categories

Figures 1 through 53: Statistical Linear Plots

a. Line Plots of Absolute/Scaled Concentrations

b. Bivariate Scatter Plotsc. Percent Differences Plots

Figures 1 through 7: Plots of Dioxin/Furan Concentrations Figures 8 through 15: Plots of Pesticide Concentrations

Figures 16 through 27: Plots of Polychlorinated Biphenyl (PCB) Concentrations

Figures 28 through 36: Plots of Polycyclic Aromatic Hydrocarbon (PAH) Concentrations

Figure 37: Plots of Total Organic Carbon (TOC) Concentrations

Figures 38 through 47: Plots of Metal Concentrations
Figure 48: Plots of Mercury Concentrations

Figure 49: Plots of Dioxins/Furans Scaled Concentrations
Figure 50: Plots of Pesticides Scaled Concentrations
Figure 51: Plots of PCBs Scaled Concentrations
Figure 52: Plots of PAHs Scaled Concentrations

Figure 53: Plots of Metals (excluding mercury) Scaled Concentrations

References

Helsel, D.L. 2005. Nondetects and Data Analysis, Statistics for Censored Environmental Data, Wiley-Interscience. 250p.

O'Brien, P.C. and T.R. Fleming. 1987. A paired Prentice-Wilcoxon test for censored paired data, Biometrics. 43, 169-180.

Table 1 Lower Passaic River 2012 Background Sediment Sampling Comparison Summary by Individual Compounds

			T			n value		1
Parameter	Number of Split Sample Pairs	Number of Split Sample Paris with Detected Concentrations	Average Ratio of CPG to EPA (for detected pairs)(1)	Percent Difference (for detected pairs) ⁽²⁾	Wilcoxon Signed Rank test ⁽³⁾	p -value Paired Prentice Wilcoxon test (4)	Statistical Difference (Yes or No) ⁽⁵⁾	Overall Split Sample Comparison (Same or Different) ⁽⁶⁾
Dioxins/Furans								
1,2,3,4,6,7,8-HpCDD	4	3	1.57±0.77	33% Outside Criteria	0.789	0.564	No	Different
1,2,3,4,6,7,8-HpCDF	4	3	1.85±0.70	67% Outside Criteria	0.181	0.083	No	Different
2,3,7,8-TCDD	4	3	0.91±0.29	Within Range	0.423	0.206	No	Same
2,3,7,8-TCDF	4	3	1.16±0.27	Within Range	1.000	0.564	No	Same
OCDD	4	3	1.90±0.70	67% Outside Criteria	0.181	0.132	No	Different
OCDF	4	3	2.38±1.17	67% Outside Criteria	0.181	0.083	No	Different
Total TCDD	4	4	1.90±0.83	25% Outside Criteria	0.100	0.058	No	Different
Pesticides								
4,4'-DDD	4	3	0.90±0.36	33% Outside Criteria	0.423	0.564	No	Same
4,4'-DDE	4	4	0.89±0.14	Within Range	0.201	0.527	No	Same
4,4'-DDT	4	3	1.05±0.38	Within Range	0.789	0.459	No	Same
Dieldrin	4	3	0.85±0.04	Within Range	0.181	0.046	No	Same
alpha-Chlordane	4	3	1.02±0.22	Within Range	0.789	1.000	No	Same
2,4'-DDD	4	3	0.59±0.61	67% Outside Criteria	0.423	0.299	No	Different
2,4'-DDE	4	2	0.84±0.05	Within Range	0.371	0.157	No	Same
2,4'-DDT	4	1	NA	Within Range	NA	0.194	No	Same
Polychlorinated Biphenyls (PCBs)								
3,3',4,4'-Tetrachlorobiphenyl (PCB 77)	4	3	1.48±0.60	33% Outside Criteria	0.181	0.083	No	Different
3,4,4',5-Tetrachlorobiphenyl (PCB 81)	4	2	1.69±1.30	50% Outside Criteria	1.000	0.527	No	Different
2,3,3',4,4'-Pentachlorobiphenyl (PCB 105)	4	3	1.22±0.52	33% Outside Criteria	0.789	0.739	No	Same
2,3,4,4',5-Pentachlorobiphenyl (PCB 114)	4	3	1.33±0.59	33% Outside Criteria	0.789	0.467	No	Different
2,3',4,4',5-Pentachlorobiphenyl (PCB 118)	4	4	1.22±0.40	25% Outside Criteria	0.584	1.000	No	Same
2',3,4,4',5-Pentachlorobiphenyl (PCB 123)	4	2	1.14±0.24	Within Range	1.000	1.000	No	Same
3,3',4,4',5-Pentachlorobiphenyl (PCB 126)	4	2	1.09±0.48	Within Range	1.000	1.000	No	Same
2,3,3',4,4',5-Hexachlorobiphenyl+2,3,3',4,4',5'-	4	3	1.10±0.30	Within Range	1.000	0.564	No	Same
Hexachlorobiphenyl (PCB 156+157)			0.05:0.44					
2,3',4,4',5,5'-Hexachlorobiphenyl (PCB 167)	4	3	0.95±0.11	Within Range	0.789	0.564	No	Same
3,3',4,4',5,5'-Hexachlorobiphenyl (PCB 169)	4	0	NA O 74 : 0 04	NA	NA 0.374	0.317	No	NA C
2,3,3',4,4',5,5'-Heptachlorobiphenyl (PCB 189)	4	2	0.71±0.21	50% Outside Criteria	0.371	0.196	No	Same
Total PCB	4	4	0.85±0.53	25% Outside Criteria	0.584	0.450	No	Same
Polycyclic Aromatic Hydrocarbons (PAHs)	4	2	1 02 10 27	Within Pango	0.789	0.564	No	Compa
Anthracene	4	3	1.02±0.37	Within Range Within Range	1.000	0.564	No	Same
Benzo[a]anthracene	4	3	1.11±0.29		0.789	0.564	No	Same
Benzo[a]pyrene	4	3	0.97±0.30	Within Range			No	Same
Chrysene	4	3	0.89±0.26	Within Range	0.789	0.564	No	Same
Fluoranthene	4	3	1.33±0.50	33% Outside Criteria	1.000	0.564	No	Different
Indeno[1,2,3-cd]pyrene Naphthalene	4	3	1.04±0.34 0.88±0.45	Within Range 33% Outside Criteria	0.789	0.564 0.879	No No	Same
·				33% Outside Criteria		0.564		Same
Phenanthrene Pyrene	4	3	1.30±0.63 1.25±0.41	33% Outside Criteria	1.000	0.564	No No	Same Same
	4	3	1.25±0.41	33/0 Outside Criteria	1.000	0.304	NO	Same
Total Organic Carbon (TOC)	4	2	2 0244 22	220/ Outside Criteria	0.422	0.266	Ne	Different
TOC Metals	4	3	2.03±1.33	33% Outside Criteria	0.423	0.366	No	Different
	4	4	0.77+0.21	25% Outside Criteria	0.201	0.225	No	Sama
Arsenic Barium	4	4	0.77±0.21 1.12±0.72	50% Outside Criteria	0.201 0.855	0.225 0.564	No No	Same Same
Cadmium	4	4	8.7±13.08	50% Outside Criteria	0.855	0.364	No No	Different
Chromium	4	4	8.7±13.08 1.14±0.37	Within Range	0.584	1.000	No	Same
	4	4	1.14±0.37 1.03±0.38	25% Outside Criteria	1.000	0.739		
Copper				25% Outside Criteria 25% Outside Criteria			No No	Same
Copper	4	4	1.01±0.46		0.855	1.000	No No	Same
Iron	4	4	0.85±0.42	25% Outside Criteria 25% Outside Criteria	0.584	0.217	No No	Same
Lead	4	4	1.37±0.25	Within Range	0.100	0.058 0.670	No	Different
Nickel	4	4	1.01±0.33	Within Range Within Range	0.855	0.870	No No	Same
Zinc			1.18±0.27				No	Same
Mercury	4	3	0.90±0.21	Within Range	1.000	0.564	No	Same

Results outside acceptance criteria are bolded.

NA = not applicable EPA = United States Environmental Protection Agency CPG = Cooperating Parties Group

Notes:

- (1) Average ratio (criteria: 0.70-1.30) with standard deviation
- (2) Percent difference criteria: no more than 16% of split samples outside of 40 to -67 %D.
- (3) Wilcoxon Signed Rank test was employed at significance level (p) of 0.05 (4) Paired Prentice Wilcoxon test was employed at significance level (p) of $0.05\,$
- (5) Limited statistical power due to small sample size (less than five)
- $(6) \ \text{If there are at least two of the three criteria (average ratio, percent different and statistical difference)} \ \text{met},$ the overall split sample comparison would be labeled "same". Otherwise, it would be "different".

Abbreviations:

1,2,3,4,6,7,8-HpCDD = 1,2,3,4,6,7,8- heptachlorodibenzo-p-dioxin 1,2,3,4,6,7,8-HpCDF = 1,2,3,4,6,7,8- heptachlorodibenzofuran 2,3,7,8-TCDD = 2,3,7,8-tetrachlorodibenzo-p-dioxin 2,3,7,8-TCDF = 2,3,7,8-tetrachlorodibenzofuran

OCDD = octachlorodibenzo-p-dioxin

 ${\sf OCDF} = octachlorodibenzo furan$ Total TCDD = total tetrachlorodibenzo-p-dioxin Total PCB = total polychlorinated biphenyl

2,4'-DDD = 2,4'-dichlorodiphenyldichloroethane

2,4'-DDE = 2,4'-dichlorodiphenyldichloroethylene

2,4'-DDT = 2,4'-dichlorodiphenyltrichloroethane 4,4'-DDD = 4,4'-dichlorodiphenyldichloroethane

4,4'-DDE = 4,4'-dichlorodiphenyldichloroethylene

4,4'-DDT = 4,4'-dichlorodiphenyltrichloroethane



Table 2
Lower Passaic River 2012 Background Sediment Sampling Comparison Summary by Chemical Categories

					p -value			
Parameter	Number of Split Sample Pairs	Number of Split Sample Paris with Detected Concentrations	Average Ratio of CPG to EPA (for detected pairs)(1)	Percent Difference (for detected pairs) (2)	Wilcoxon Signed Rank test ⁽³⁾	Paired Prentice Wilcoxon test	Statistical Difference (Yes or No)	Overall Split Sample Comparison (Same or Different) ⁽⁵⁾
Dioxins/Furans	24	18	1.63±0.79	39% Outside Criteria	0.009	0.027	Yes	Different (CPG higher than EPA)
Pesticides	32	22	0.87±0.30	14% Outside Criteria	0.041	0.015	Yes	Same
Polychlorinated Biphenyls (PCBs)	48	31	1.20±0.49	19% Outside Criteria	0.207	0.652	No	Same
Polycyclic Aromatic Hydrocarbons (PAHs)	36	27	1.09±0.38	15% Outside Criteria	0.710	0.904	No	Same
Metals (excluding mercury)	40	40	1.82±4.32 ⁽⁶⁾	23% Outside Criteria	0.317	0.311	No	Same

Results outside acceptance criteria are bolded.

NA = not applicable

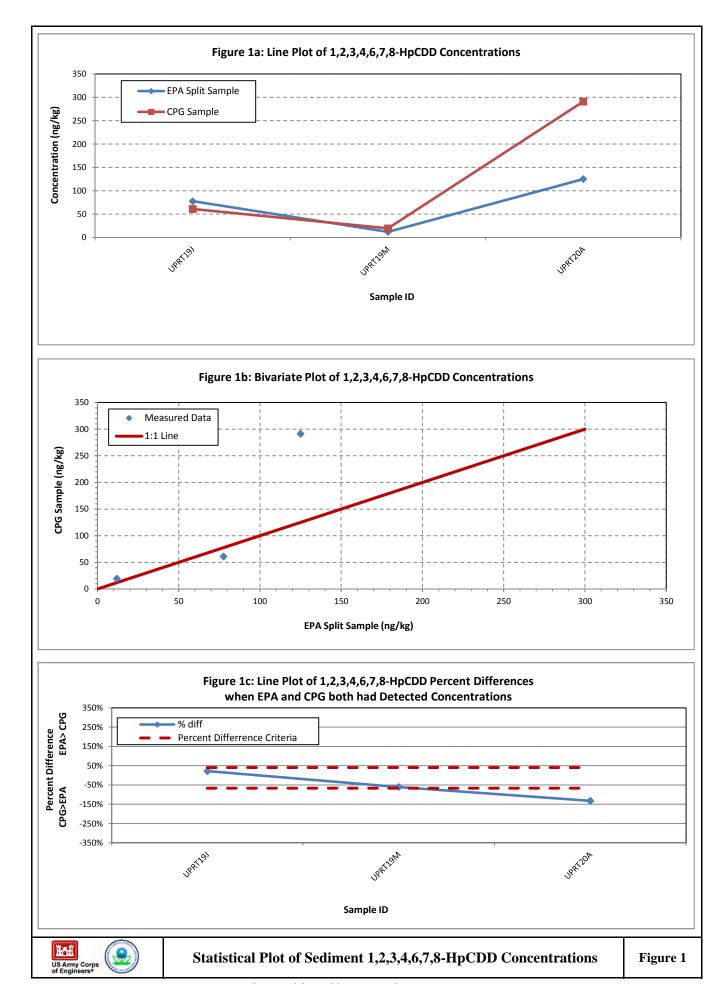
EPA = United States Environmental Protection Agency

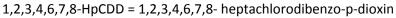
CPG = Cooperating Parties Group

Notes:

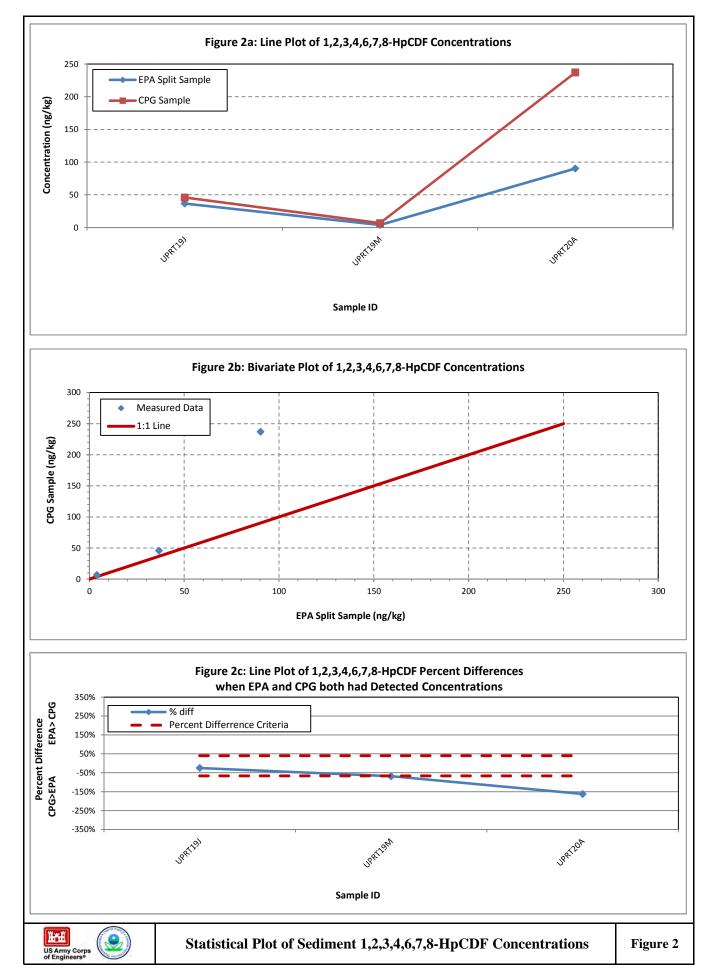
- (1) Average ratio (criteria: 0.70-1.30) with standard deviation
- (2) Percent difference criteria: no more than 16% of split samples outside of 40 to -67 %D.
- (3) Wilcoxon Signed Rank test was employed at significance level (p) of 0.05
- (4) Paired Prentice Wilcoxon test was employed at significance level (p) of 0.05
- (5) If there are at least two of the three criteria (average ratio, percent different and statistical difference) met, the overall split sample comparison would be labeled "same". Otherwise, it would be "different".
- (6) Although average ratio (1.82) is not within criterion, the actual average ratio can be within criterion because of the high standard deviation (4.32).

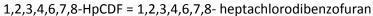




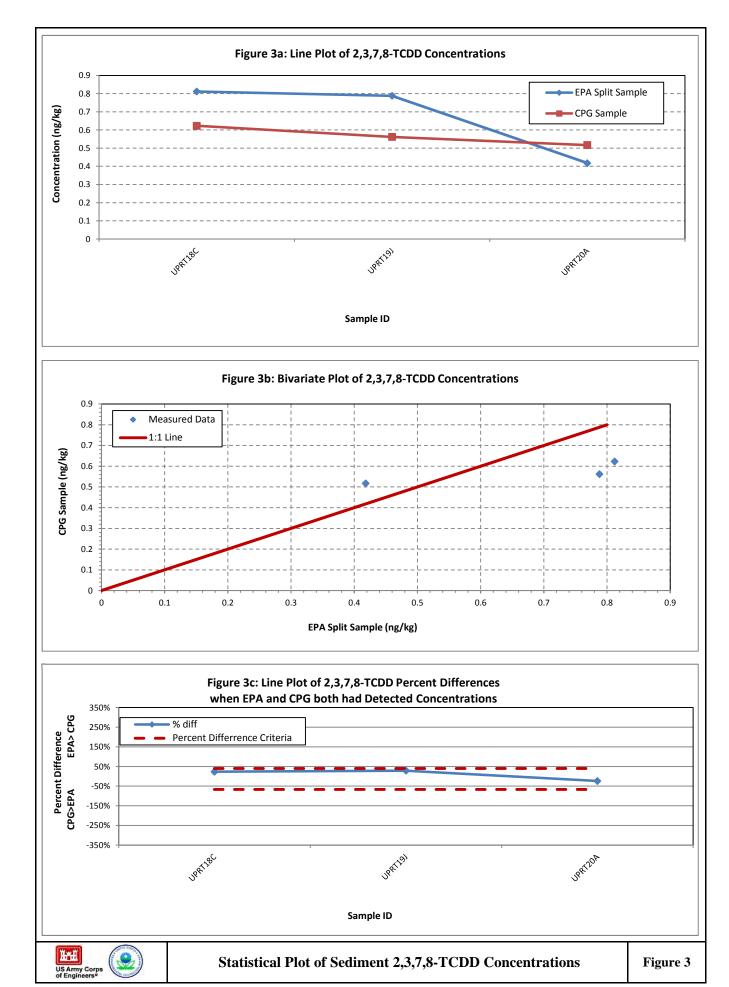


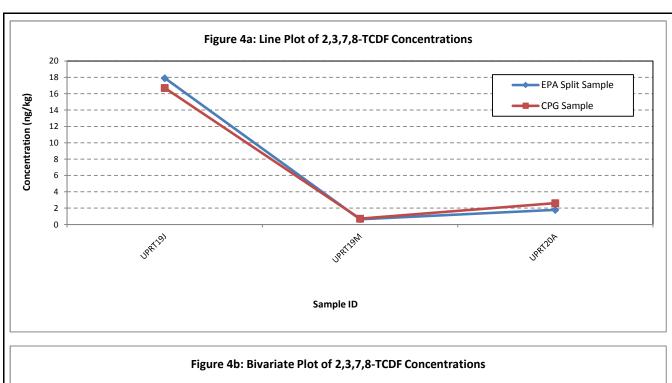


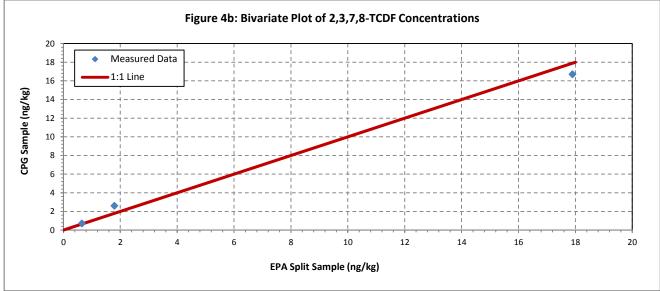


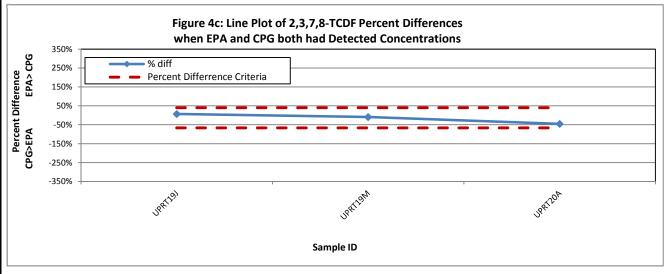






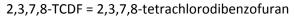




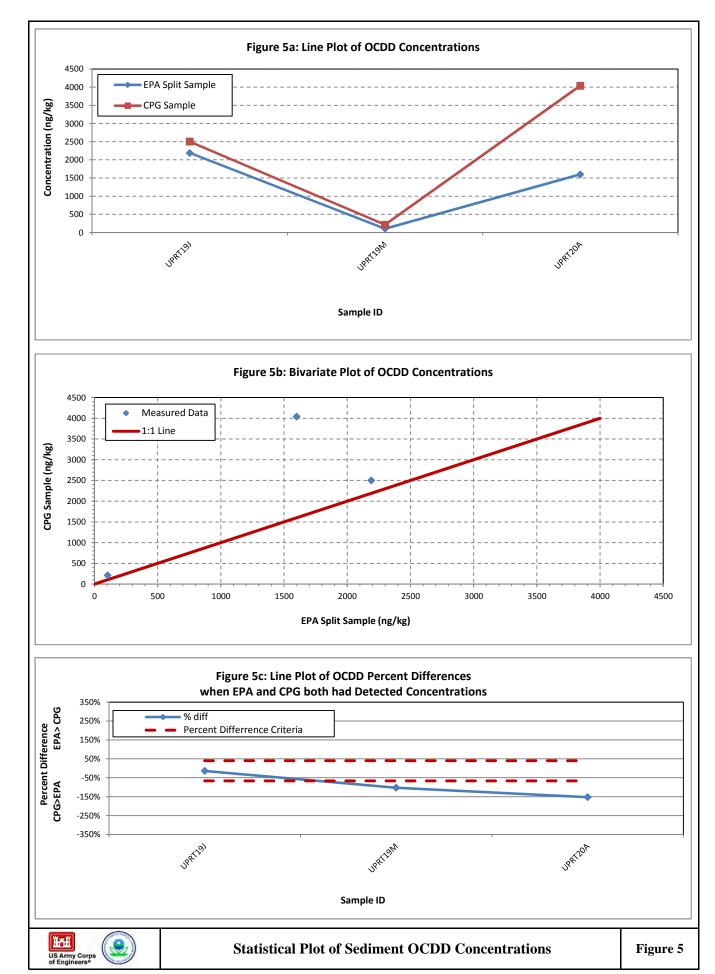




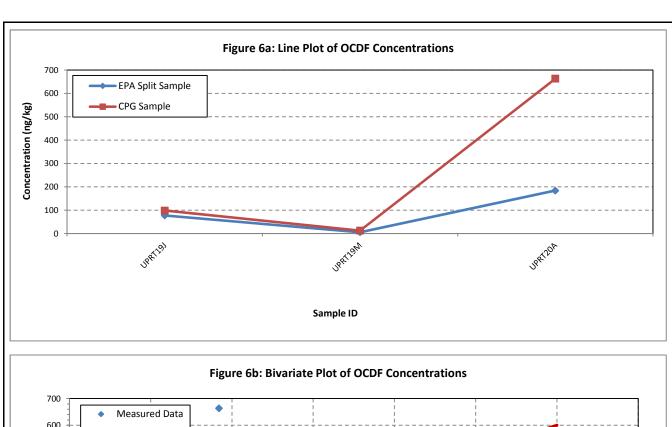
Statistical Plot of Sediment 2,3,7,8-TCDF Concentrations

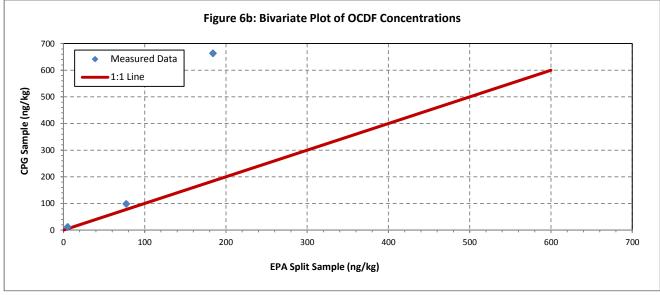


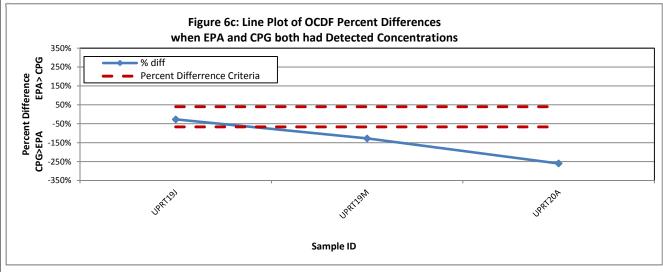










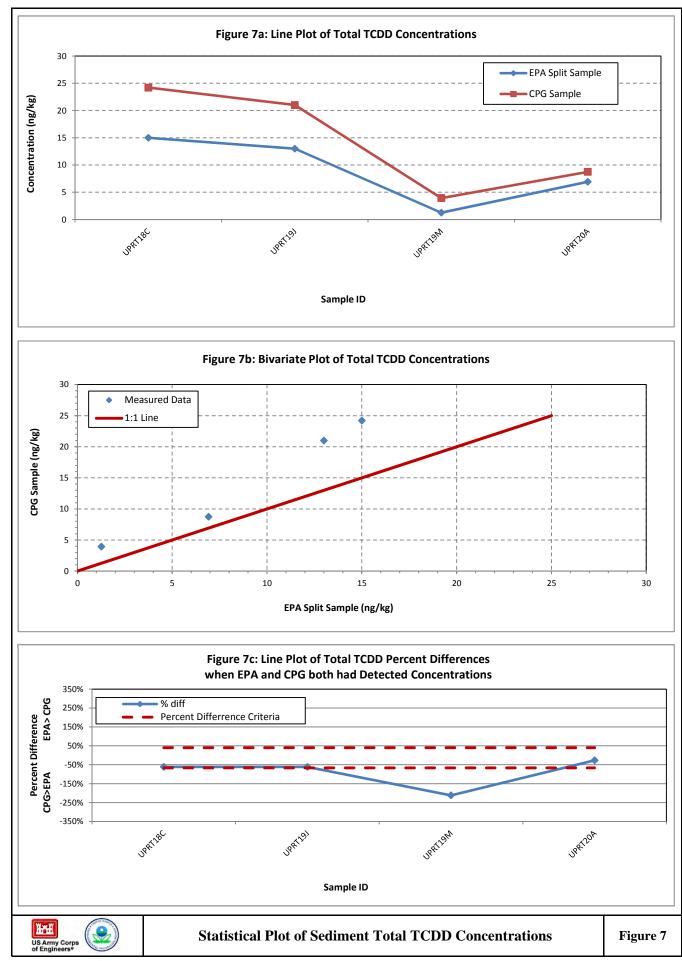


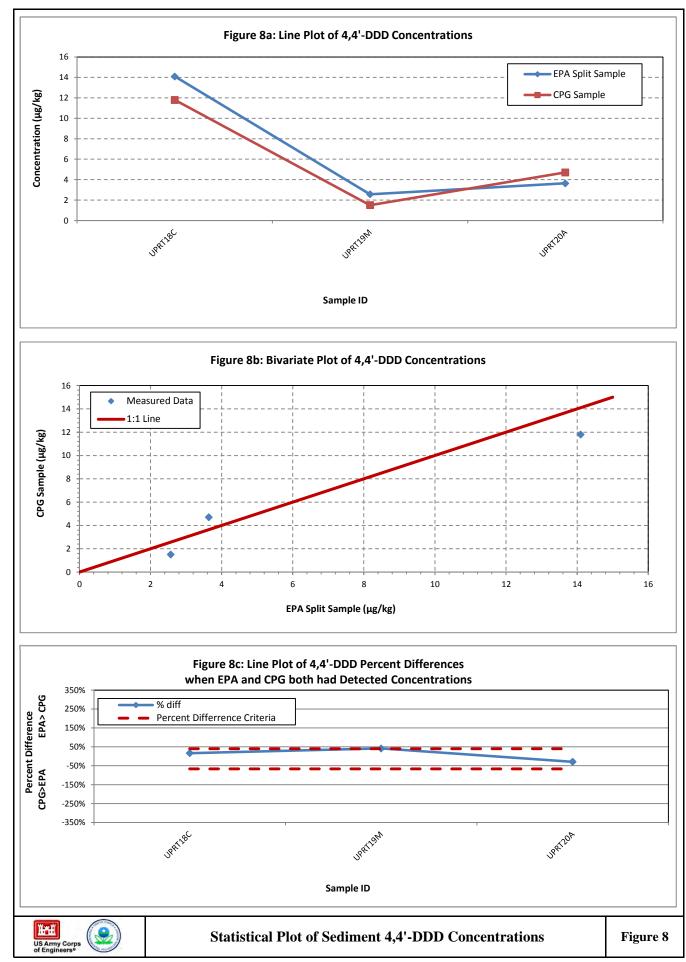


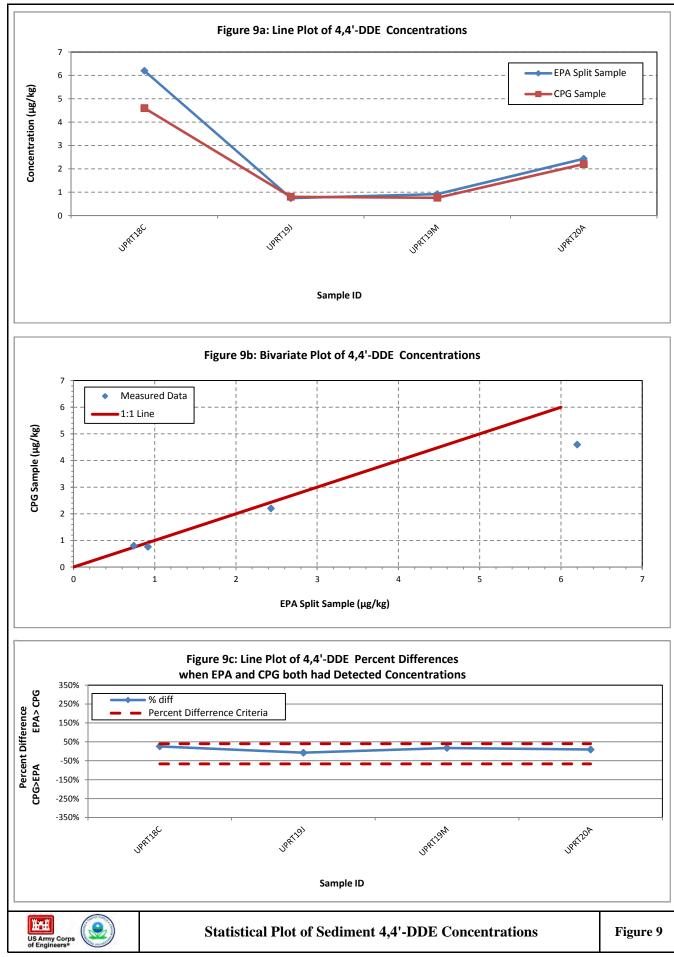
Statistical Plot of Sediment OCDF Concentrations

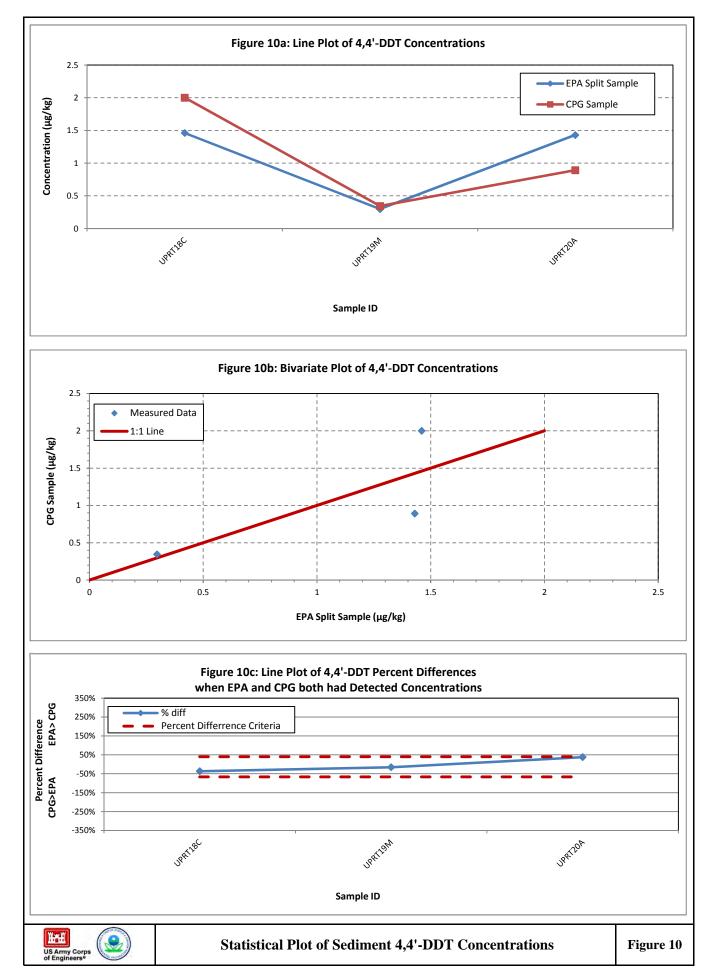






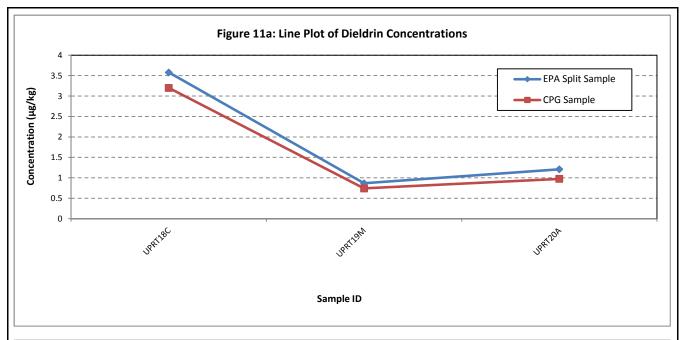


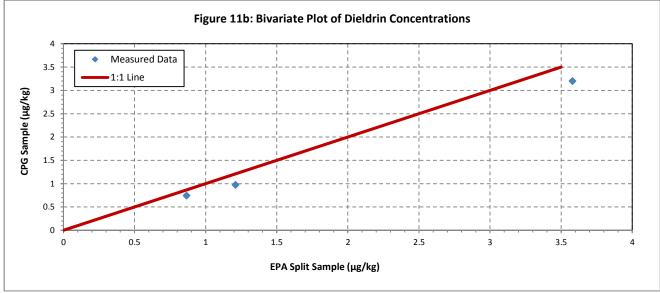


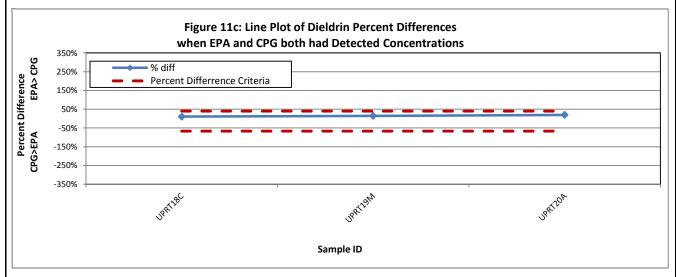










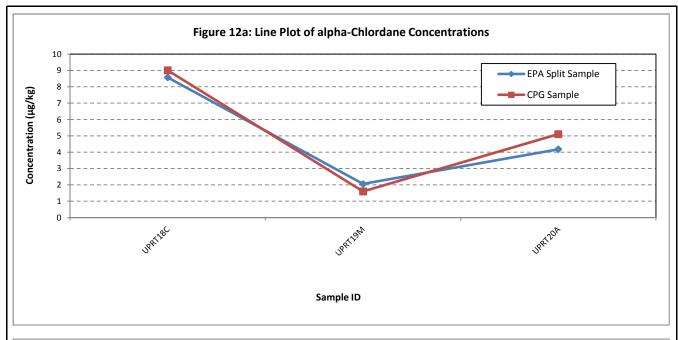


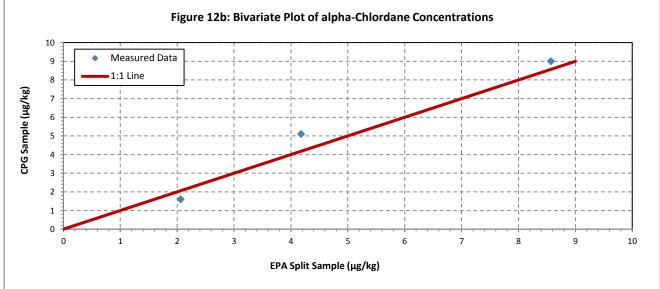


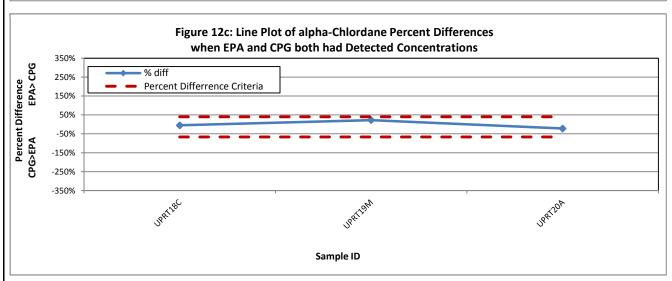


Statistical Plot of Sediment Dieldrin Concentrations







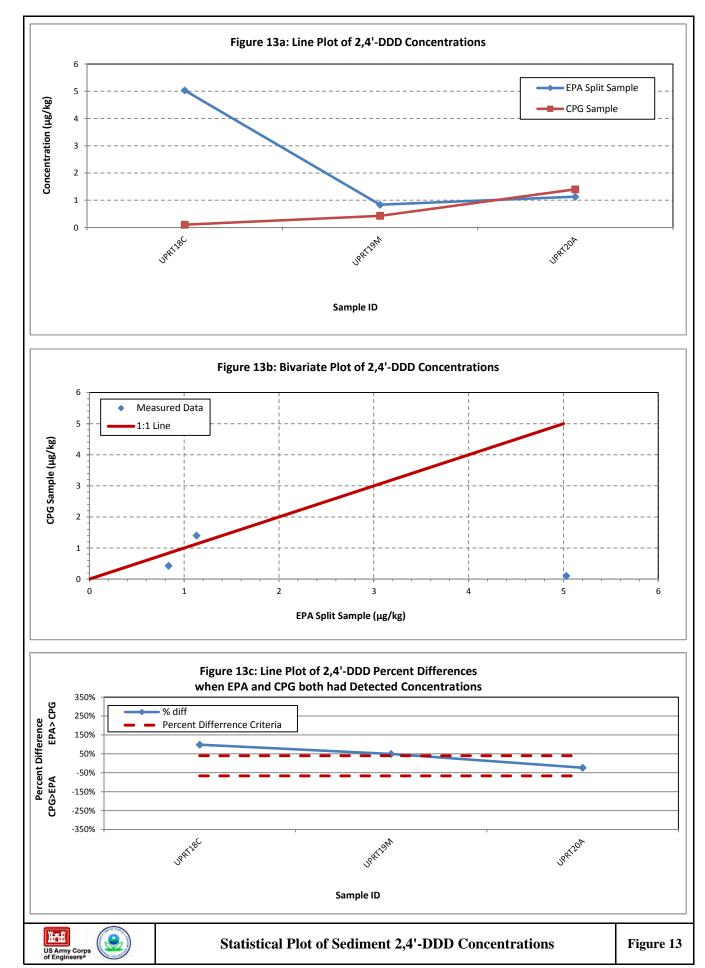




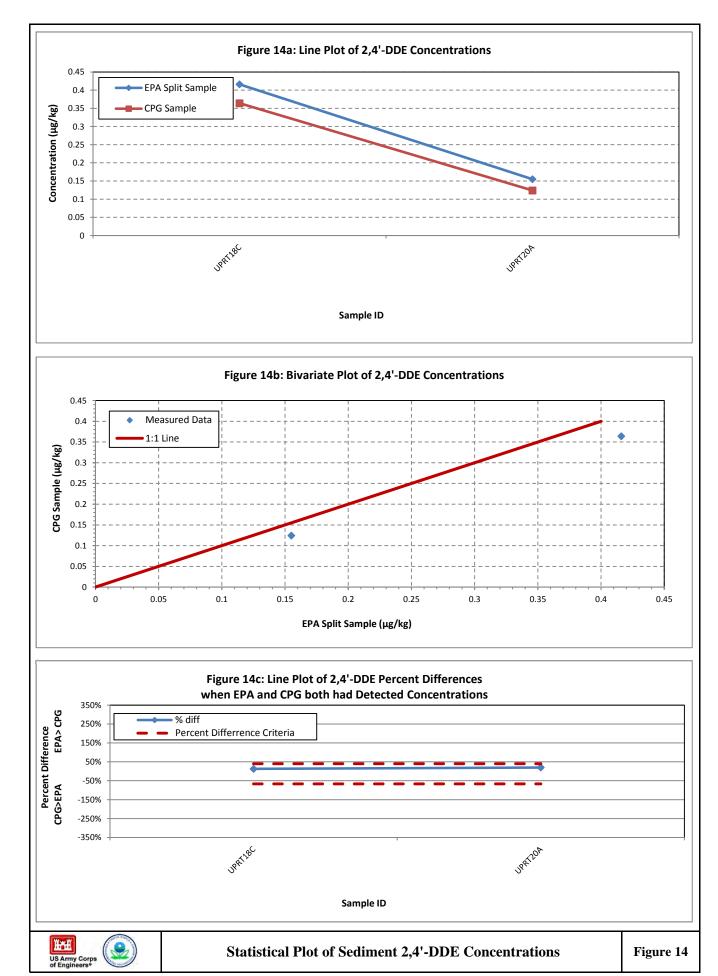


Statistical Plot of Sediment gamma-Chlordane Concentrations



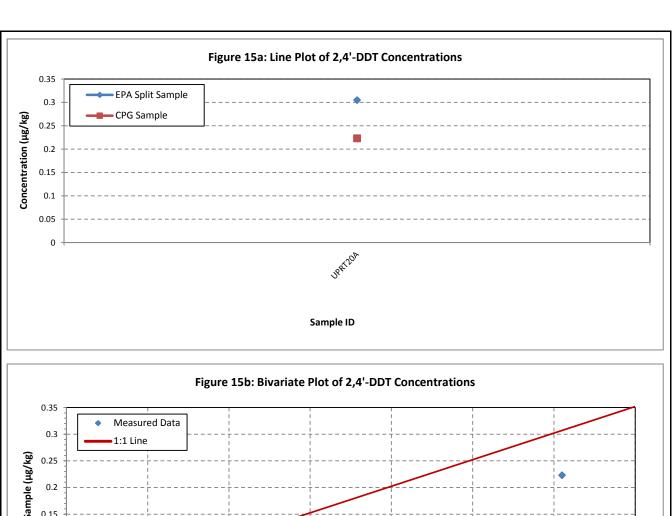


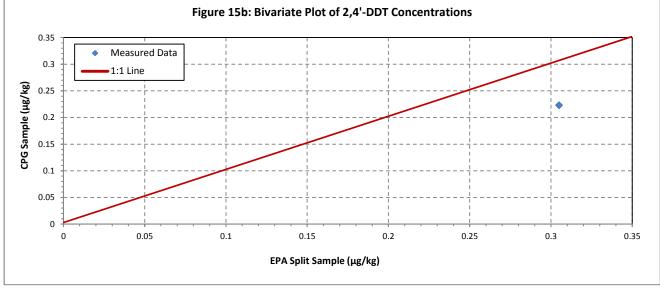


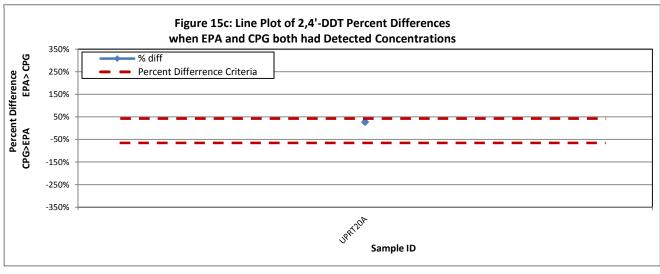






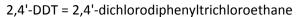




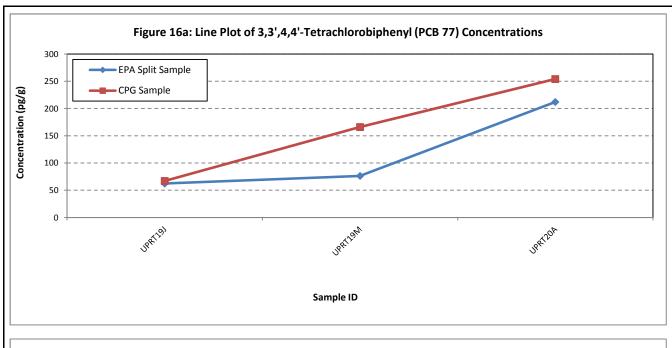


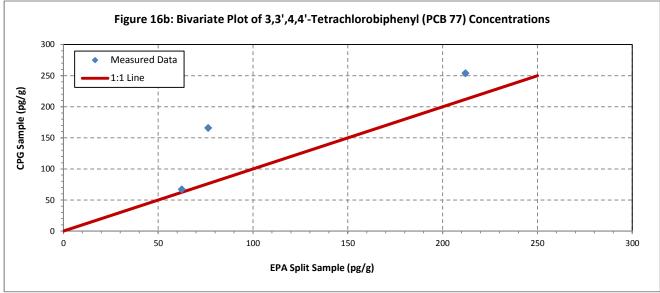


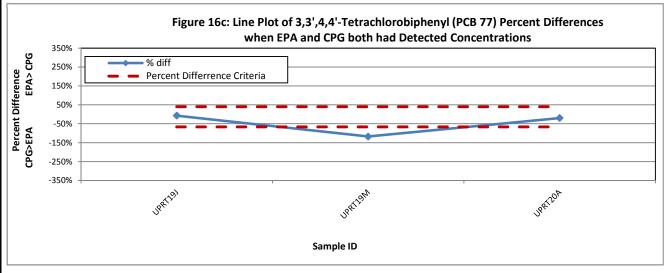
Statistical Plot of Sediment 2,4'-DDT Concentrations









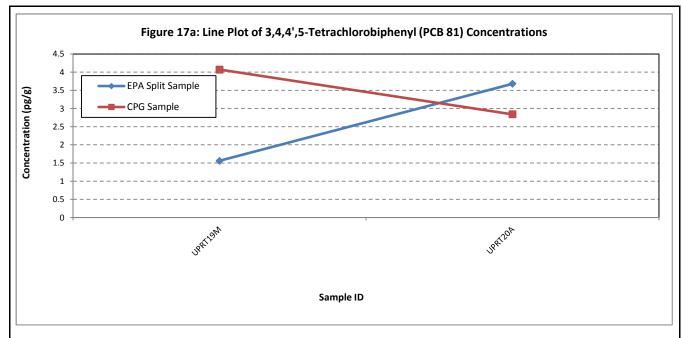


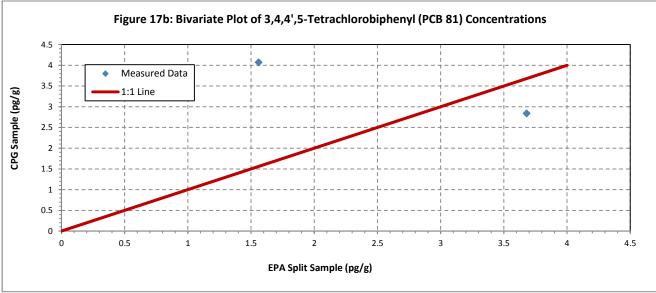


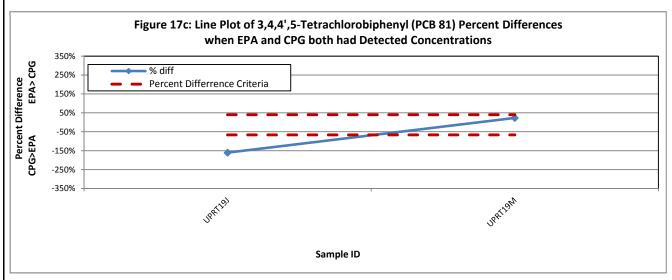
Statistical Plot of Sediment 3,3',4,4'-Tetrachlorobiphenyl (PCB 77) Concentrations









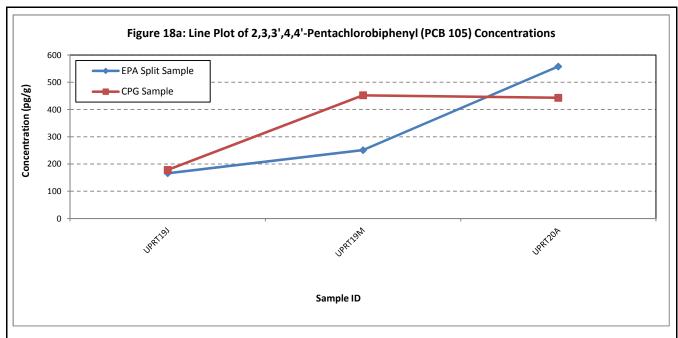


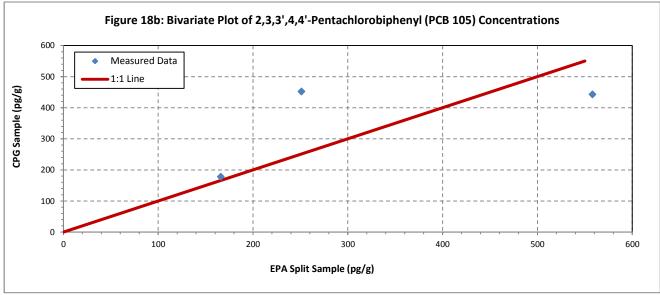


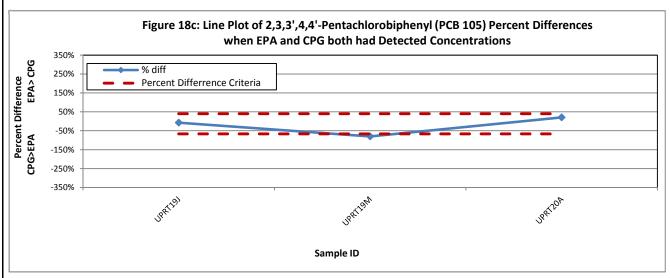
Statistical Plot of Sediment 3,4,4',5-Tetrachlorobiphenyl (PCB 81) Concentrations









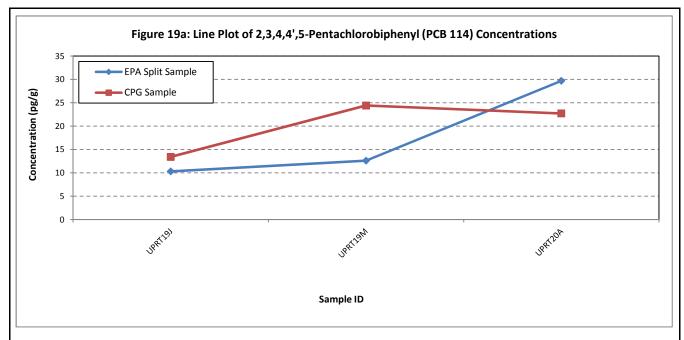


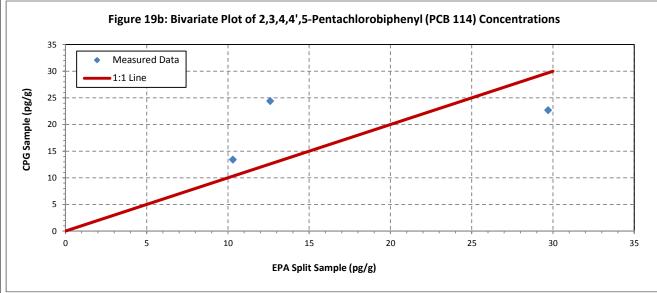


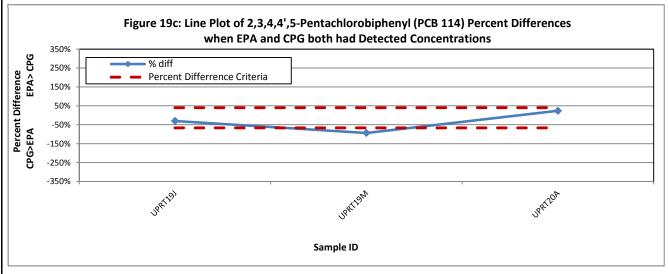
Statistical Plot of Sediment 2,3,3',4,4'-Pentachlorobiphenyl (PCB 105) Concentrations









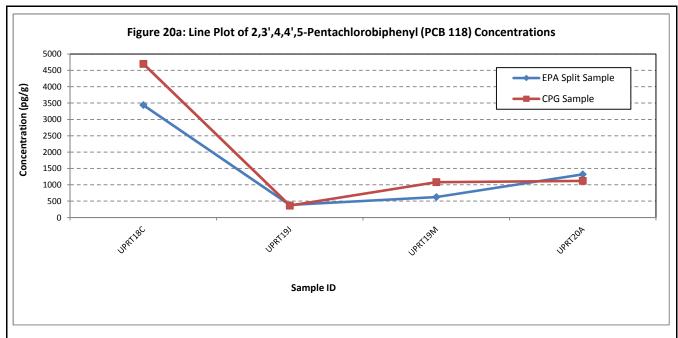


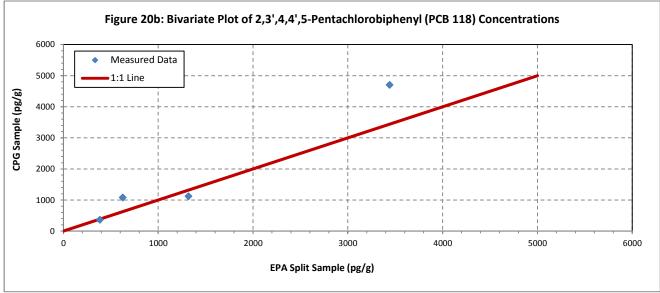


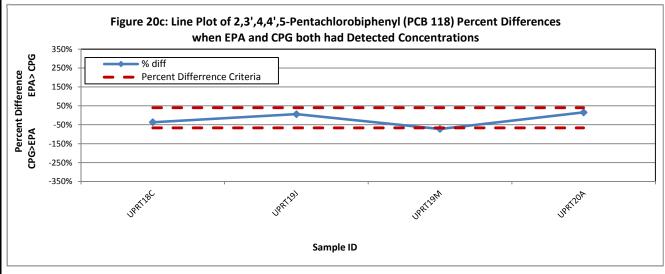
Statistical Plot of Sediment 2,3,4,4',5-Pentachlorobiphenyl (PCB 114) Concentrations









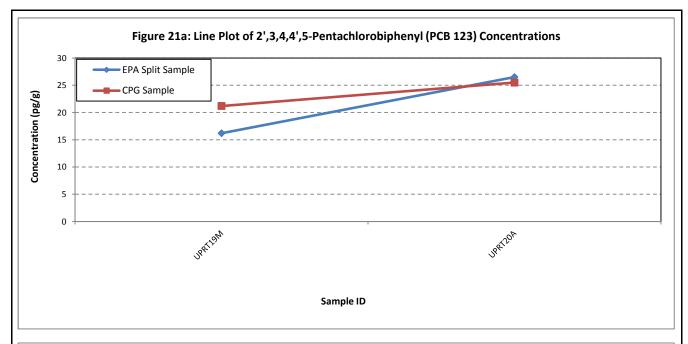


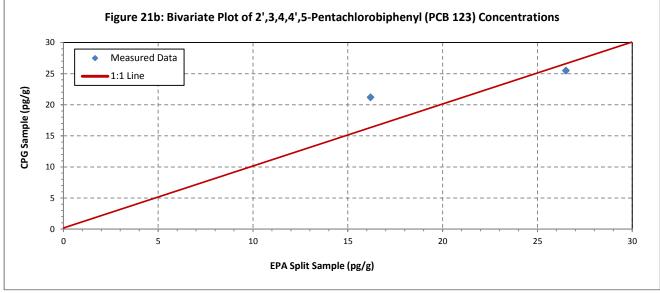


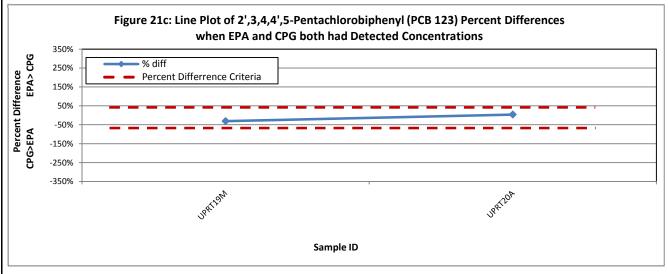
Statistical Plot of Sediment 2,3',4,4',5-Pentachlorobiphenyl (PCB 118) Concentrations









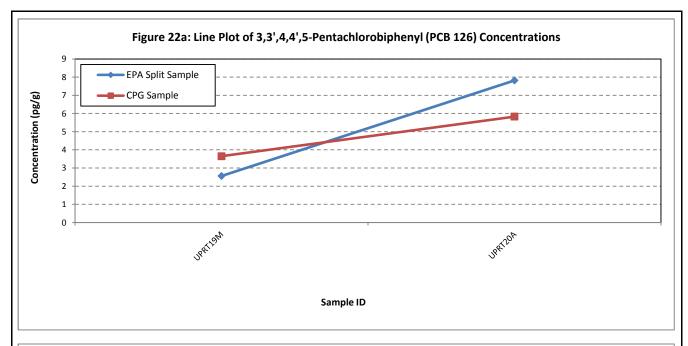


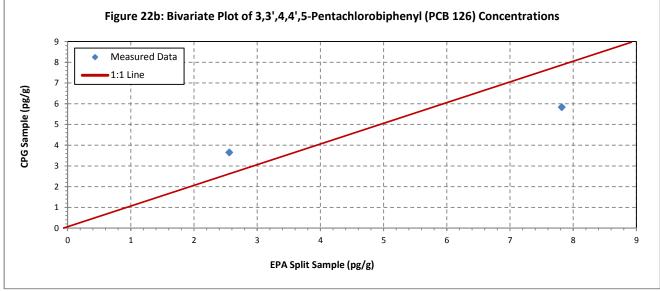


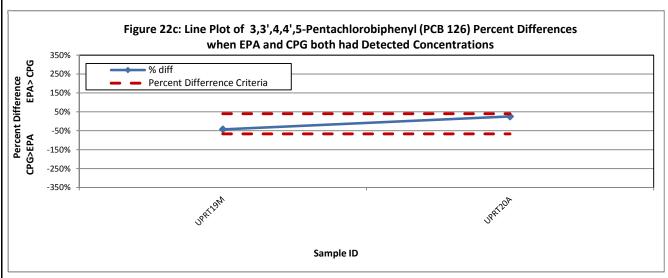
Statistical Plot of Sediment 2,3',4,4',5'-Pentachlorobiphenyl (PCB 123) Concentrations









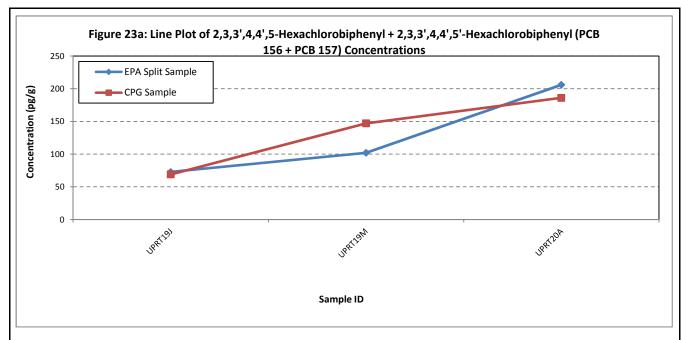


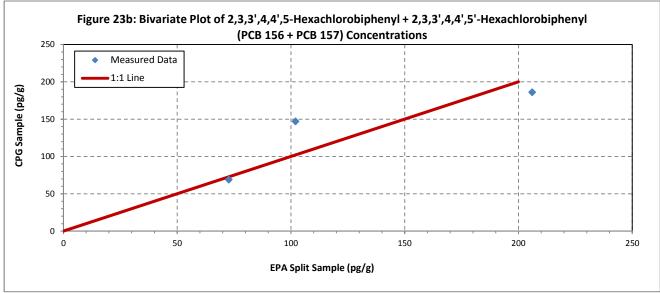


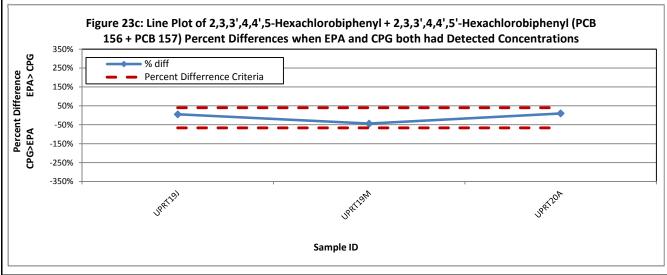
Statistical Plot of Sediment 3,3',4,4',5-Pentachlorobiphenyl (PCB 126) Concentrations









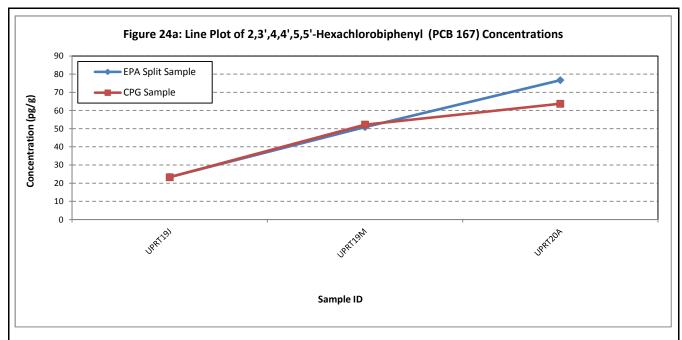


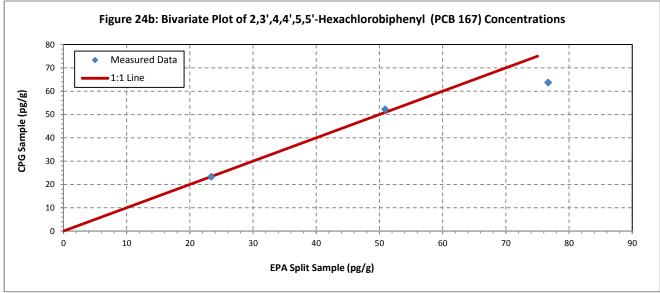


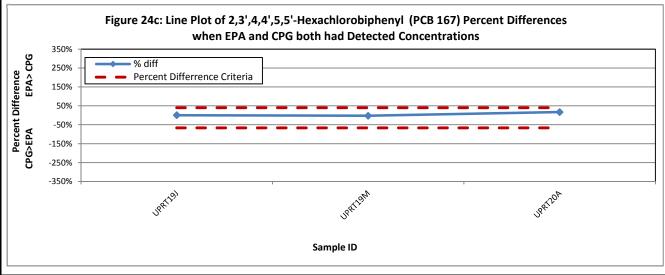
Statistical Plot of Sediment 2,3,3',4,4',5-Hexachlorobiphenyl + 2,3,3',4,4',5'-Hexachlorobiphenyl (PCB 156 + PCB 157) Concentrations













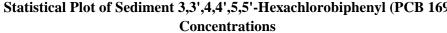
Statistical Plot of Sediment 2,3',4,4',5,5'-Hexachlorobiphenyl (PCB 167) Concentrations



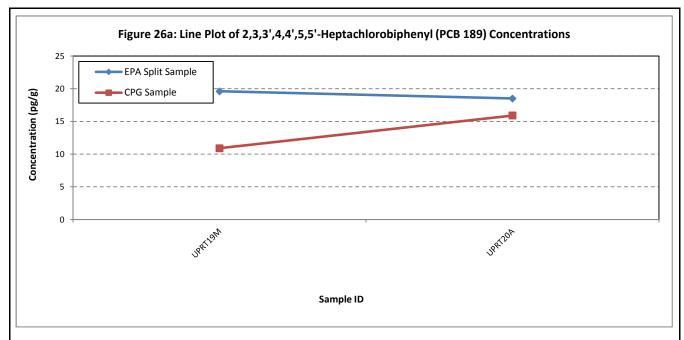


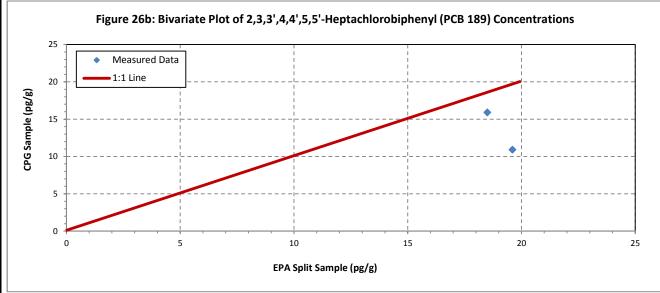
No comparison possbile bed	cause none of the sample locations had		
detected concentrations on	both USEPA and CPG samples		
	· · · · · · · · · · · · · · · · · · ·		
1	cause none of the sample locations had		
detected concentrations on	both USEPA and CPG samples		
Γ			
No comparison possbile because none of the sample locations had			
detected concentrations on both USEPA and CPG samples			
W. W.	Statistical Diet of Sediment 2 21 4 41 5 51 House Line Line 1 (DCD 160)		
US Army Corps	Statistical Plot of Sediment 3,3',4,4',5,5'-Hexachlorobiphenyl (PCB 169)	Figure 25	
US Army Corps	Concentrations	riguit 23	

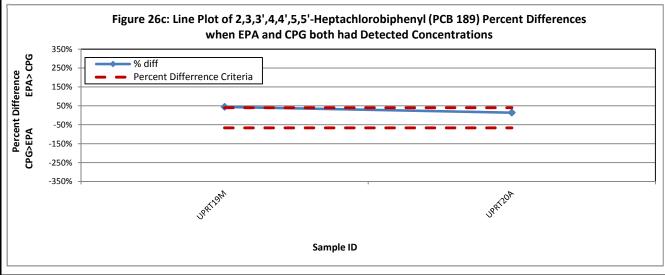










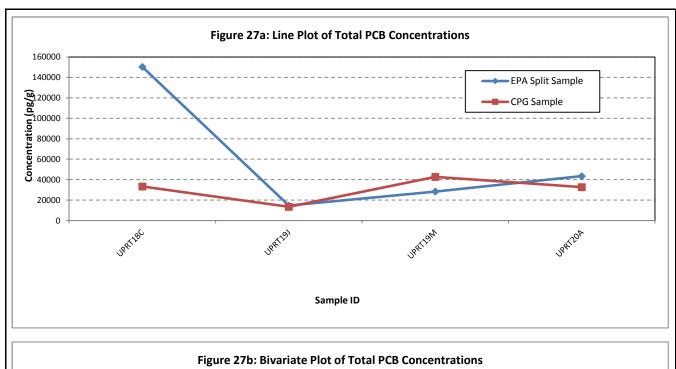


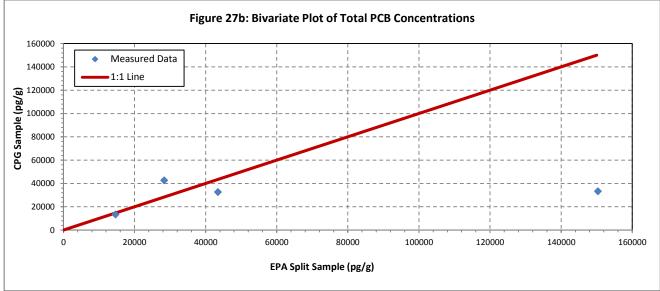


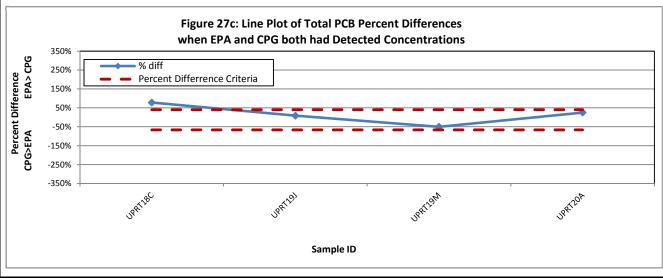
Statistical Plot of Sediment 2,3,3',4,4',5,5'-Heptachlorobiphenyl (PCB 189) Concentrations









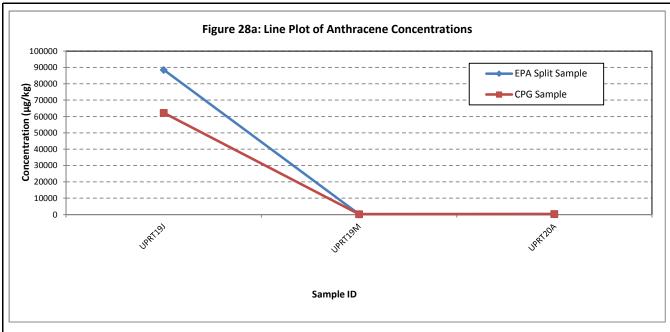


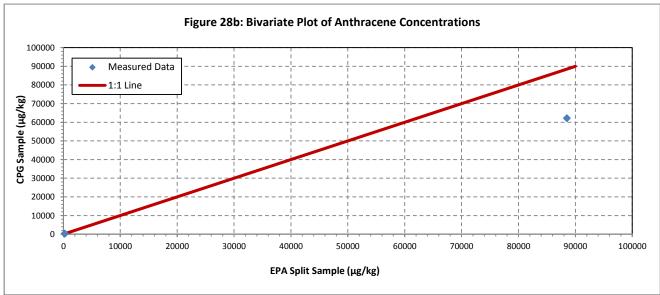


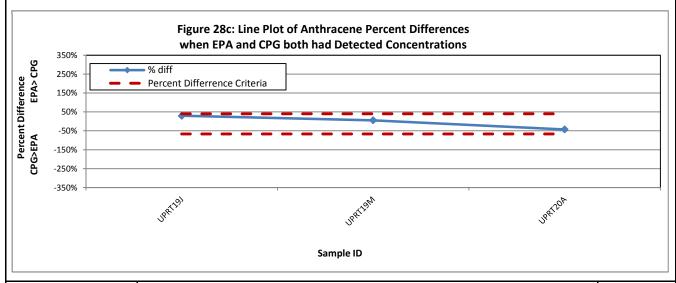
Statistical Plot of Sediment Total PCB Concentrations







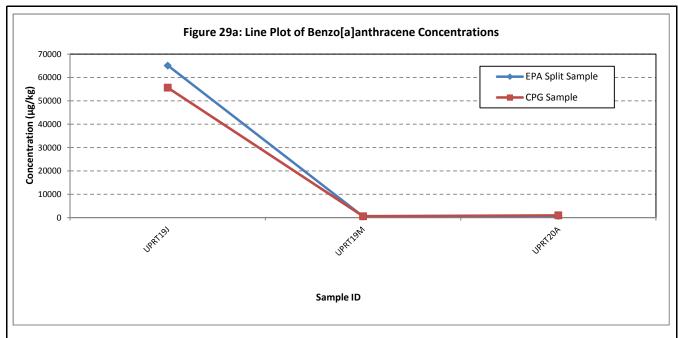


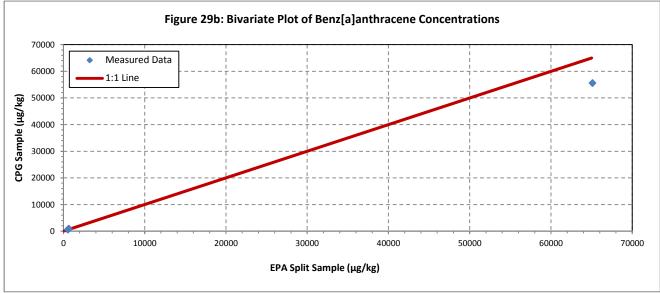


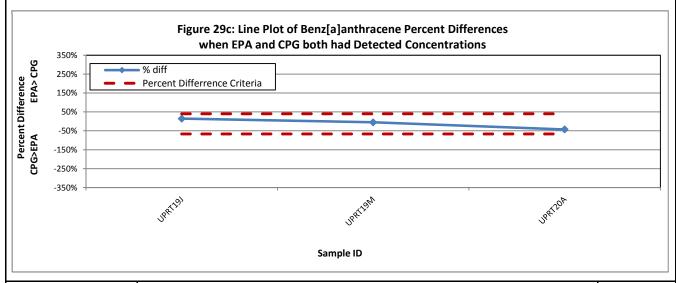


Statistical Plot of Sediment Anthracene Concentrations





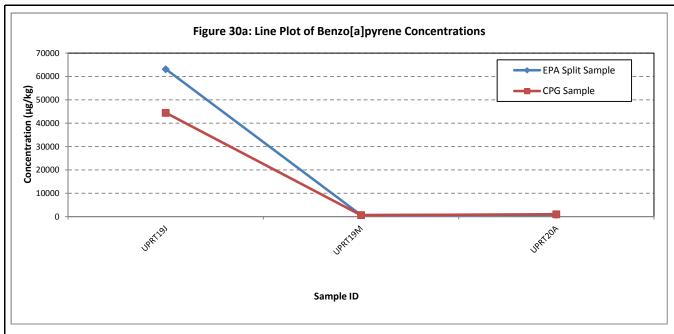


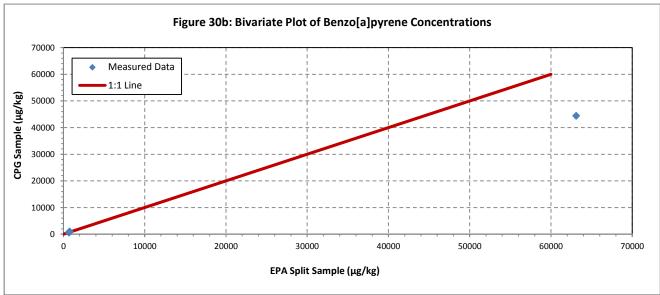


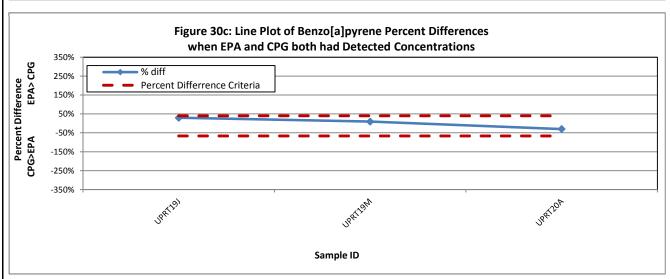


Statistical Plot of Sediment Benz[a]anthracene Concentrations







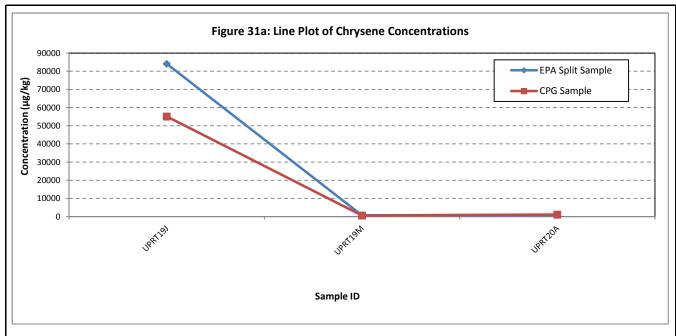


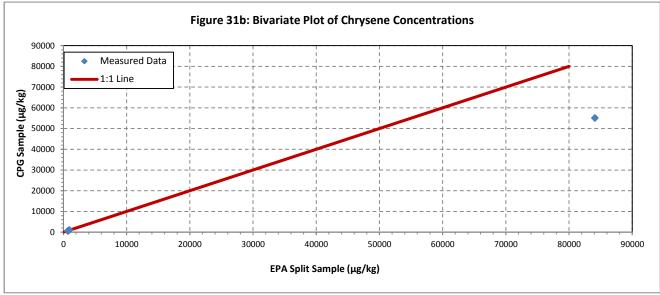


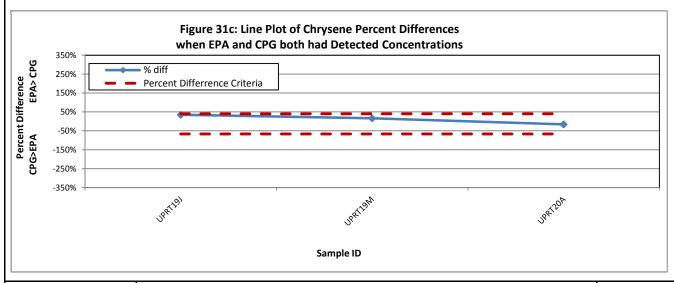


Statistical Plot of Sediment Benzo[a]pyrene Concentrations





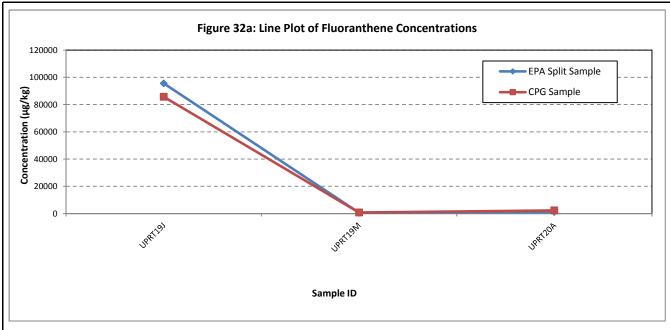


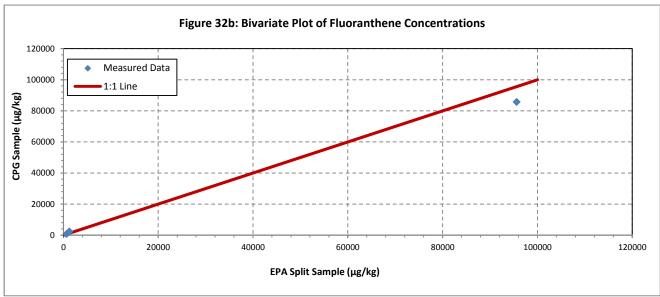


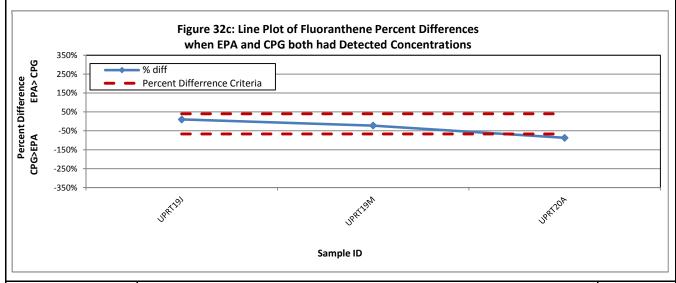


Statistical Plot of Sediment Chrysene Concentrations





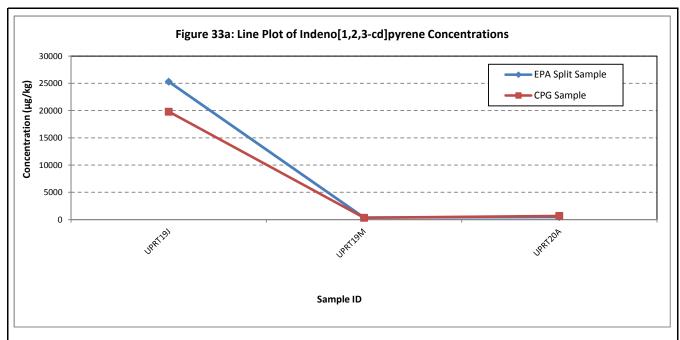


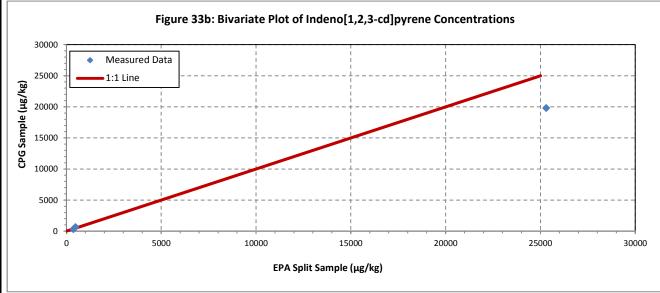


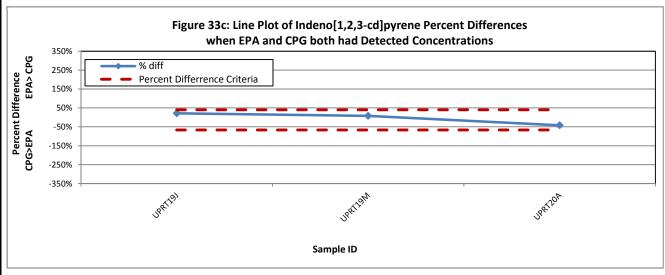


Statistical Plot of Sediment Fluoranthene Concentrations





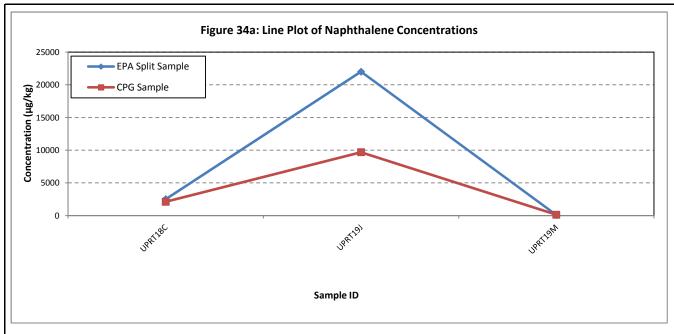


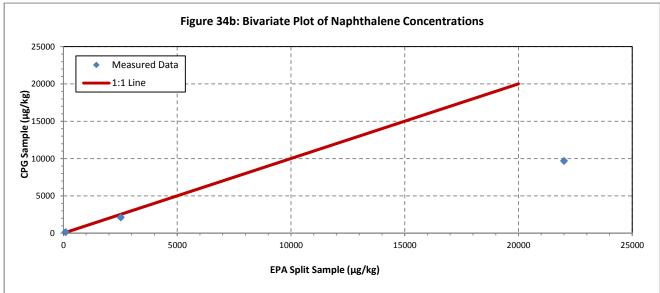


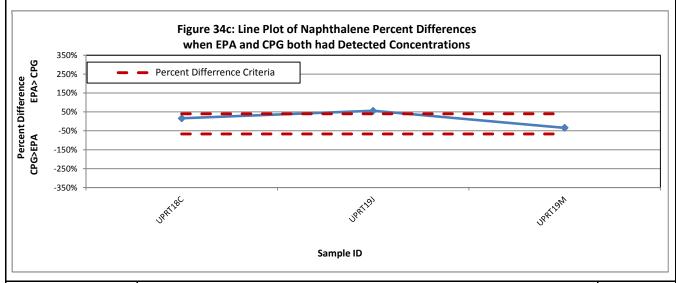


Statistical Plot of Sediment Indeno[1,2,3-cd]pyrene Concentrations





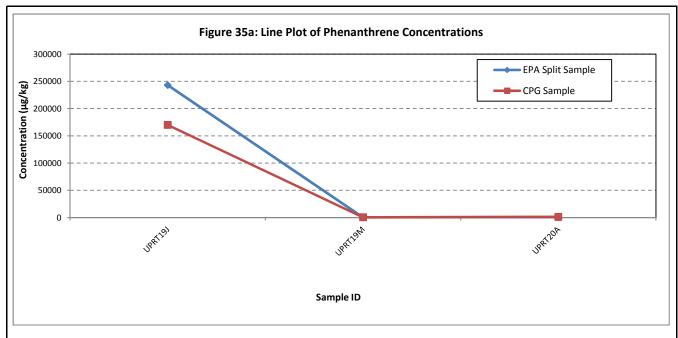


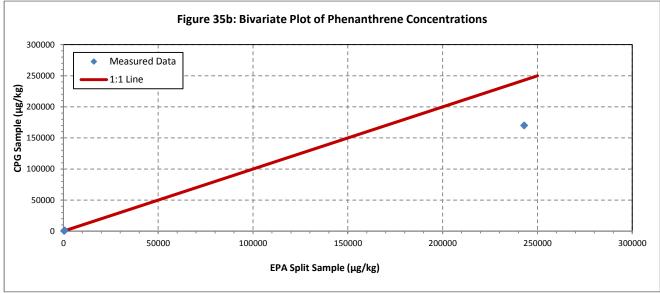


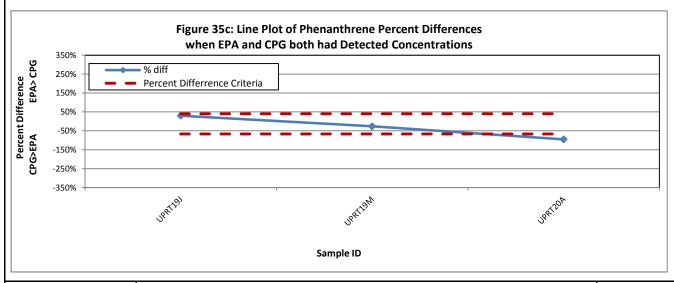


Statistical Plot of Sediment Naphthalene Concentrations





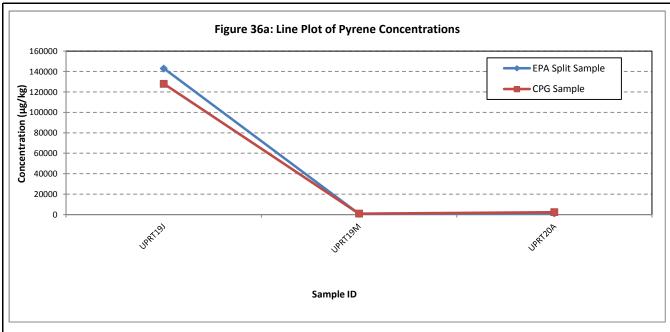


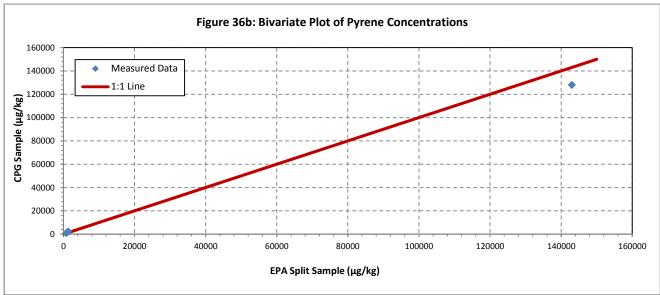


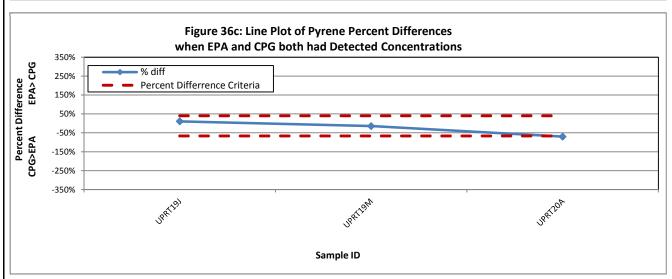


Statistical Plot of Sediment Phenanthrene Concentrations





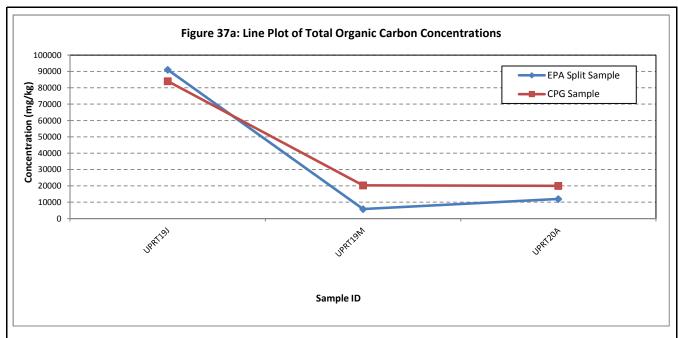


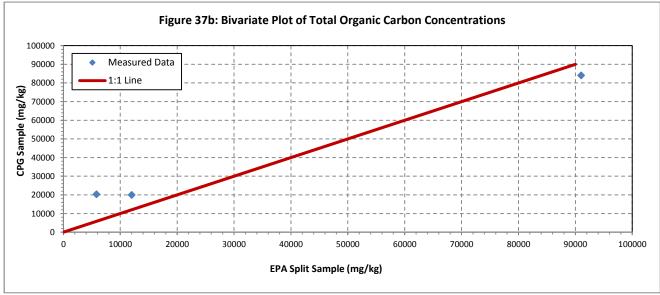


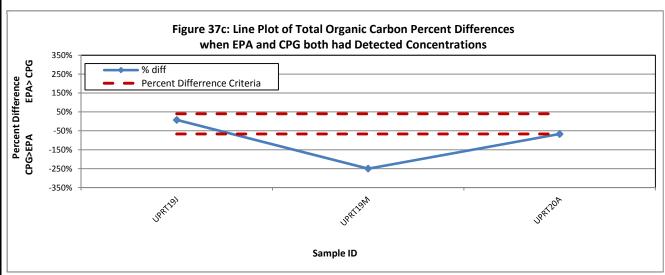


Statistical Plot of Sediment Pyrene Concentrations





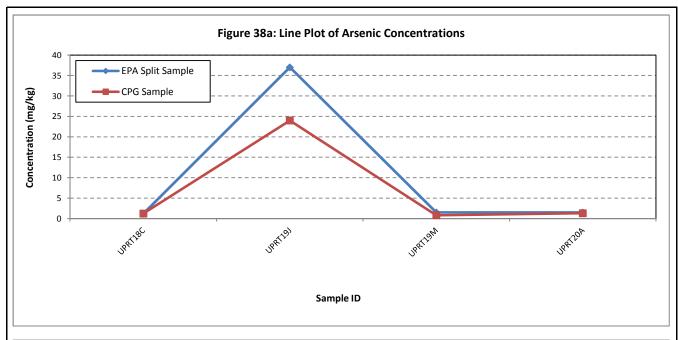


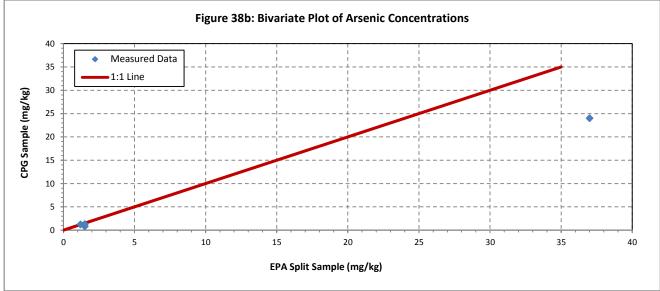


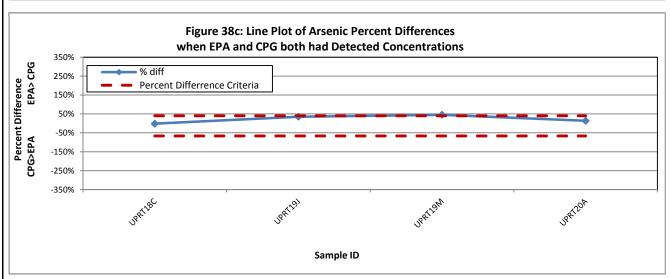


Statistical Plot of Sediment Total Organic Carbon Concentrations







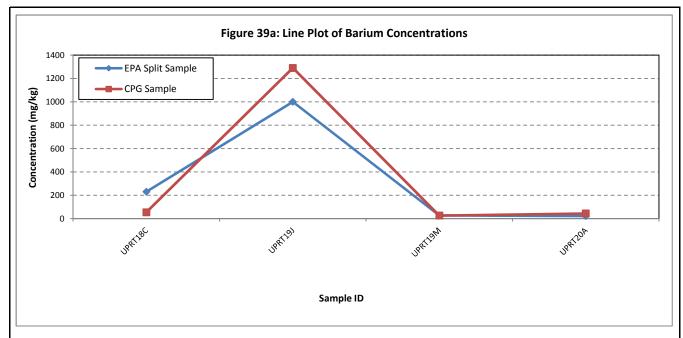


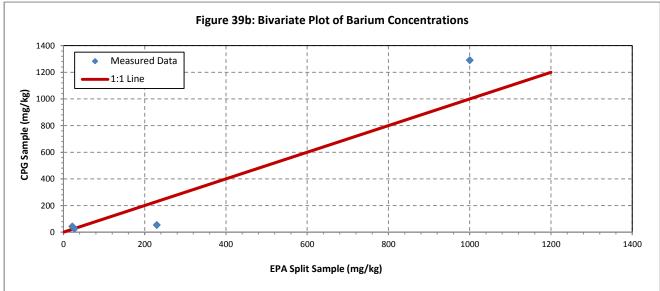


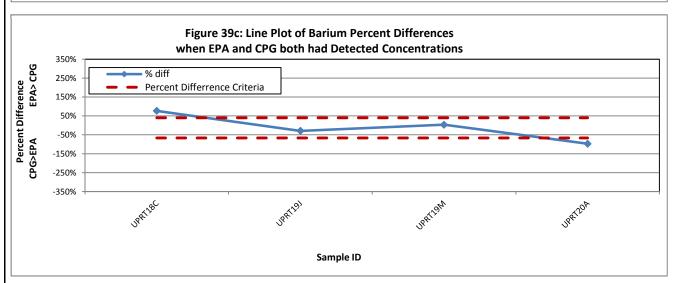


Statistical Plot of Sediment Arsenic Concentrations







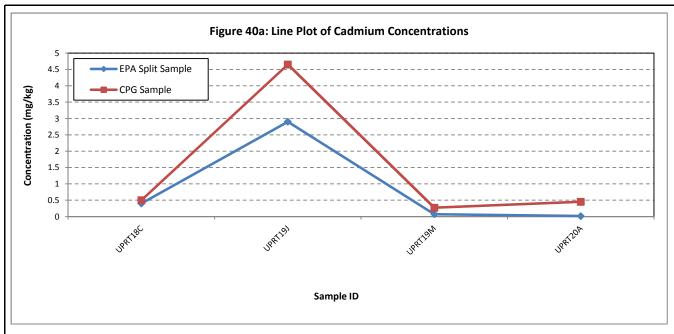


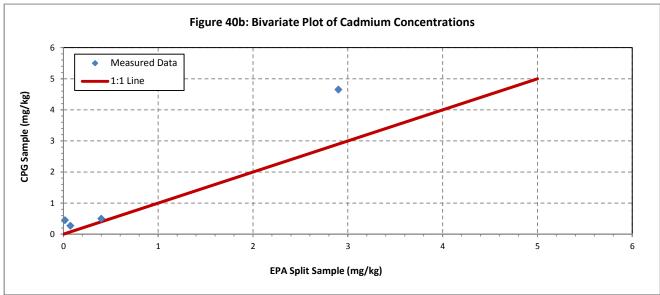


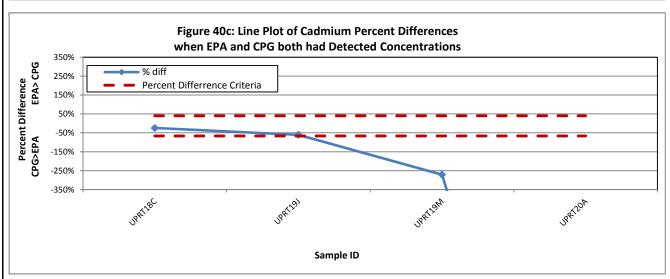


Statistical Plot of Sediment Barium Concentrations

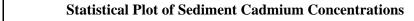




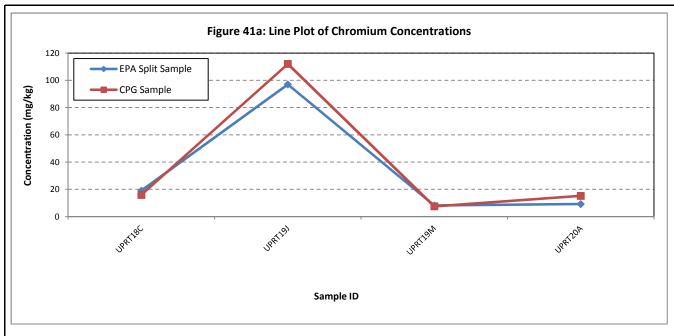


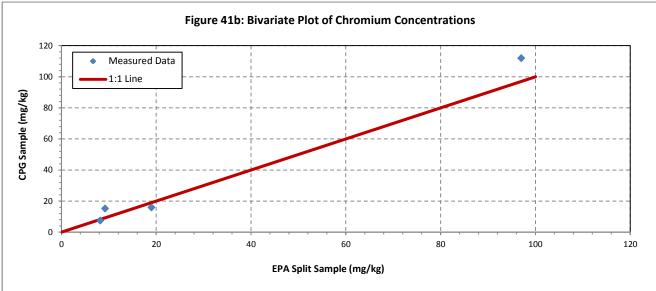


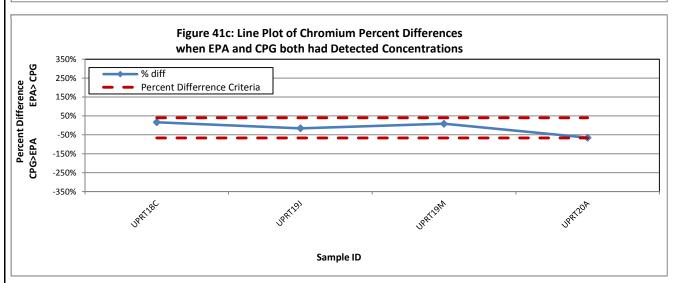










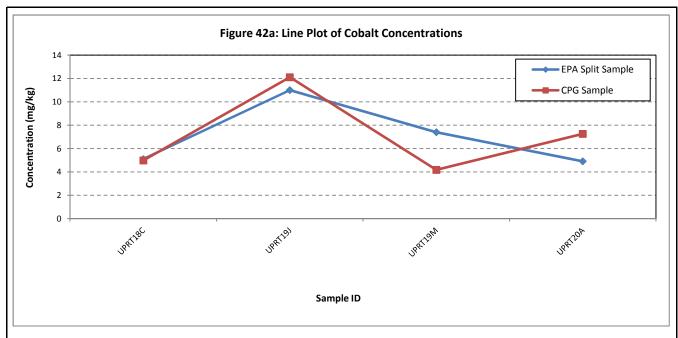


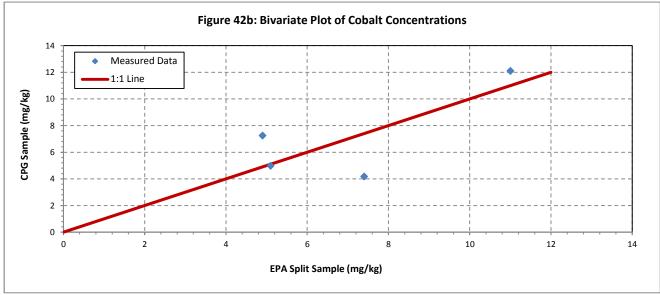


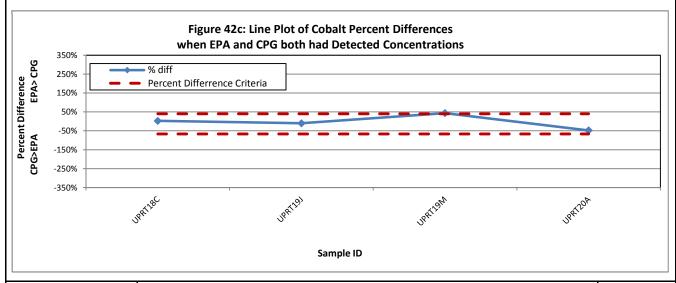


Statistical Plot of Sediment Chromium Concentrations







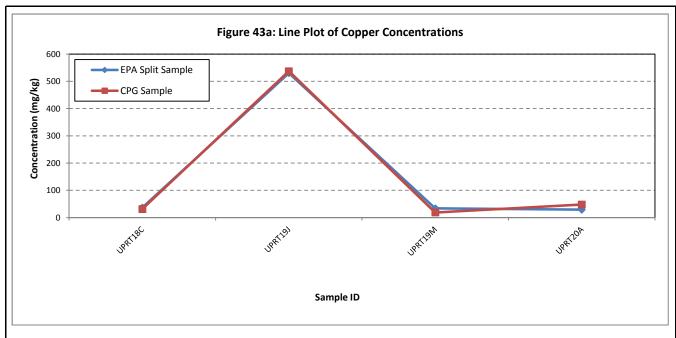


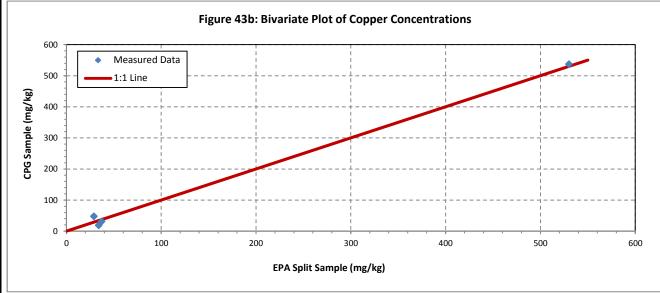


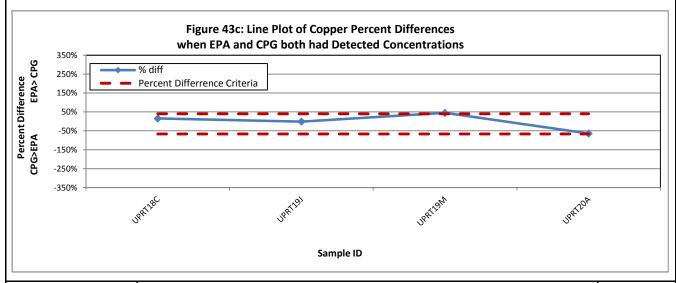


Statistical Plot of Sediment Cobalt Concentrations





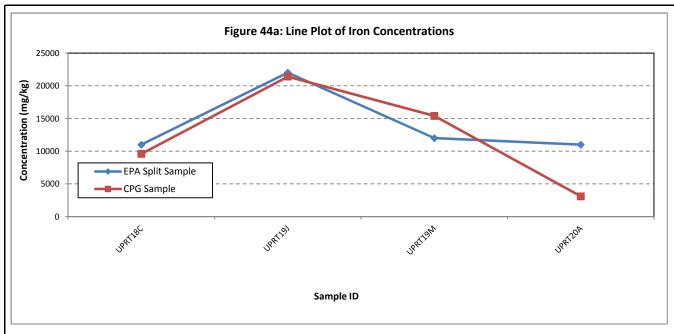


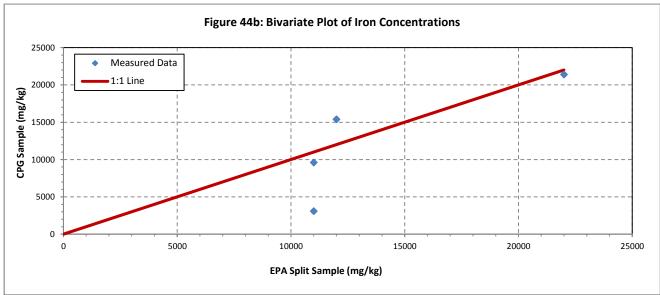


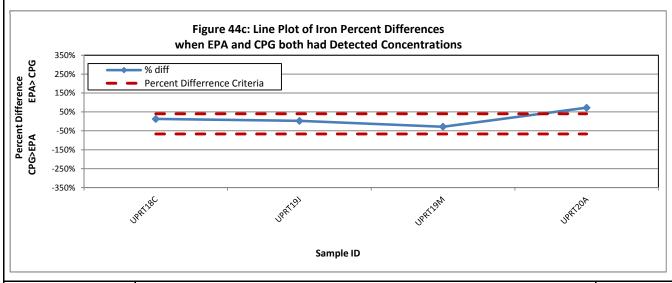


Statistical Plot of Sediment Copper Concentrations





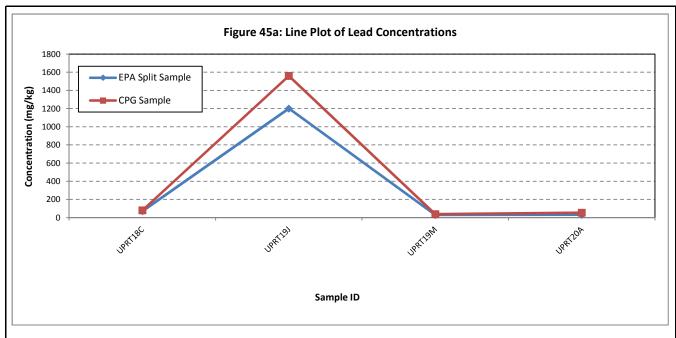


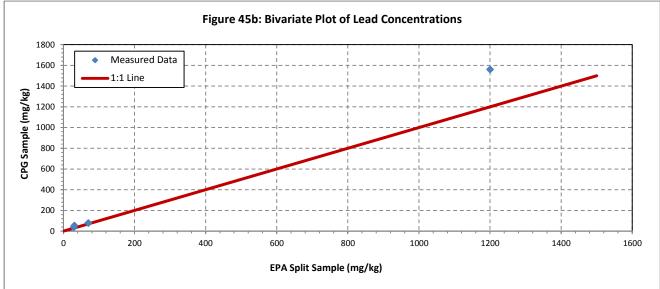


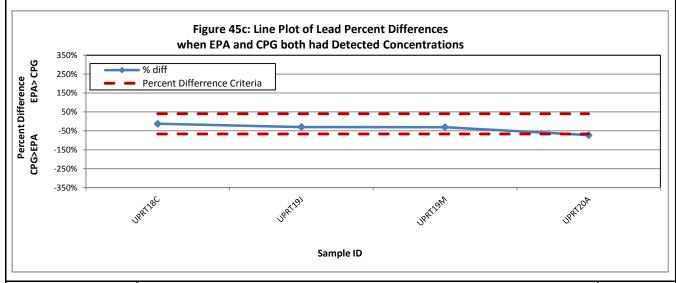


Statistical Plot of Sediment Iron Concentrations





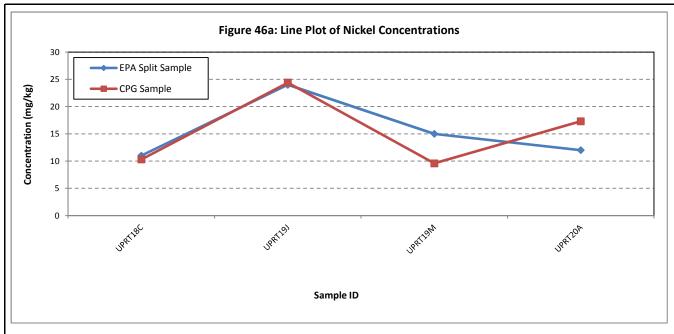


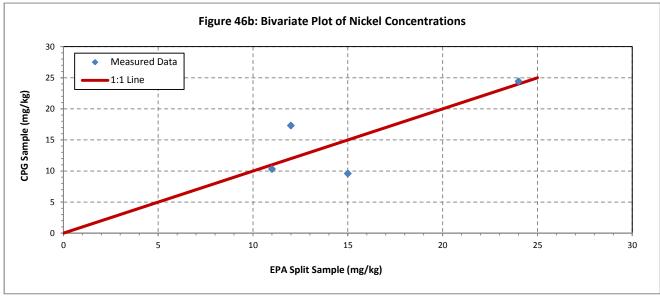


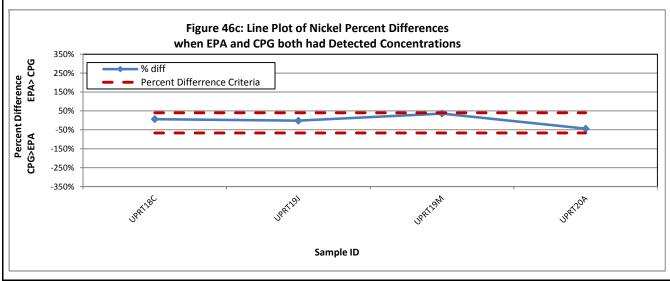


Statistical Plot of Sediment Lead Concentrations







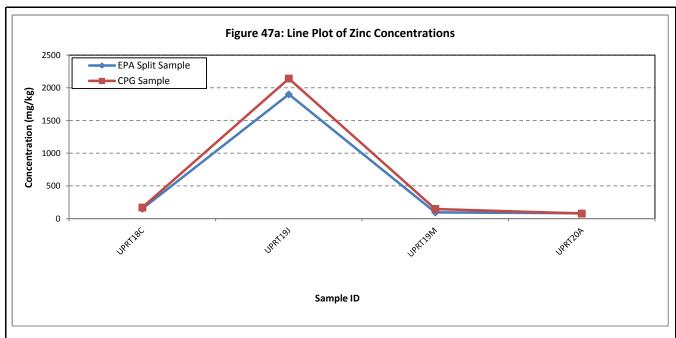


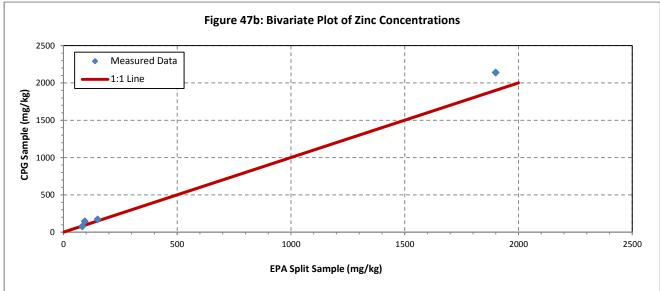


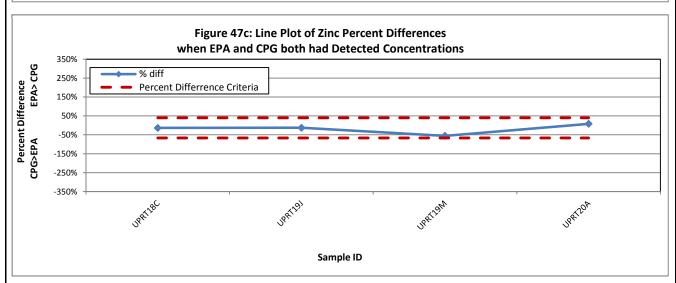


Statistical Plot of Sediment Nickel Concentrations







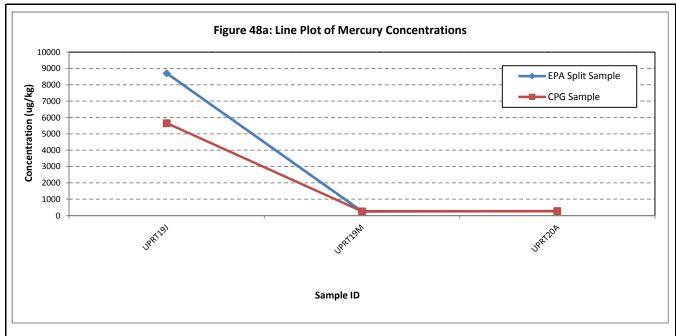


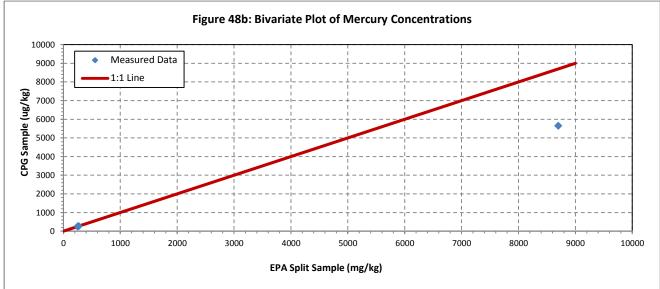


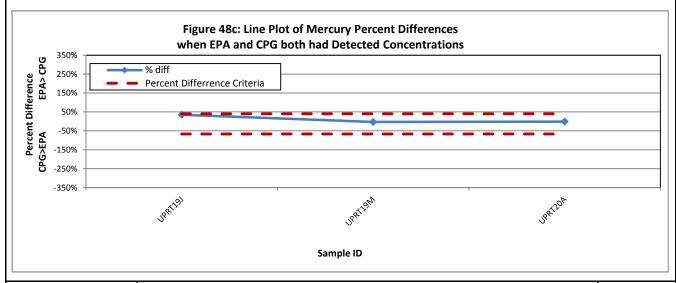


Statistical Plot of Sediment Zinc Concentrations





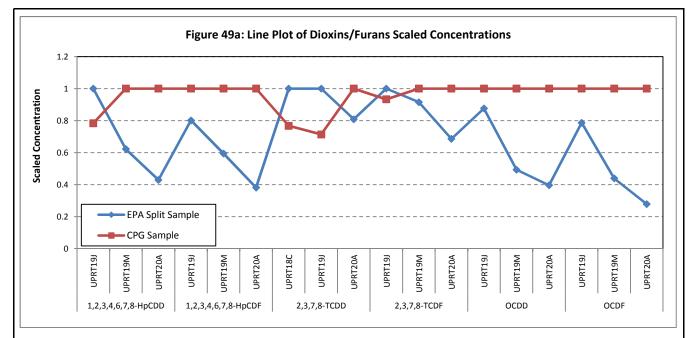


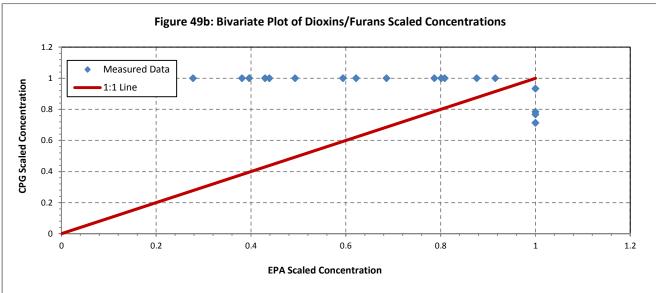


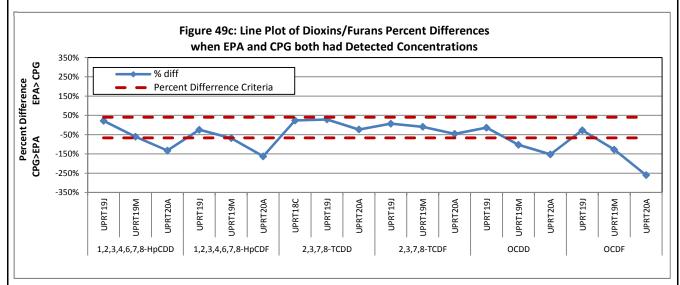


Statistical Plot of Sediment Mercury Concentrations











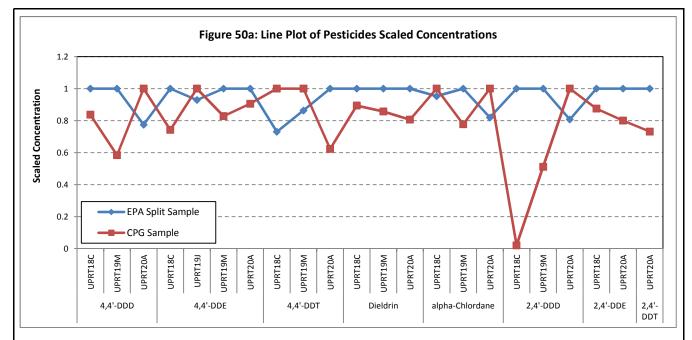
Statistical Plot of Sediment Dioxins/Furans Scaled Concentrations

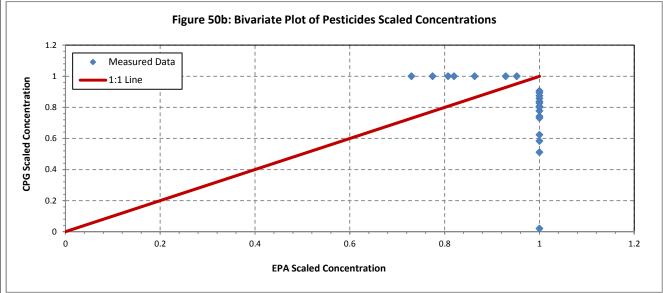
Figure 49

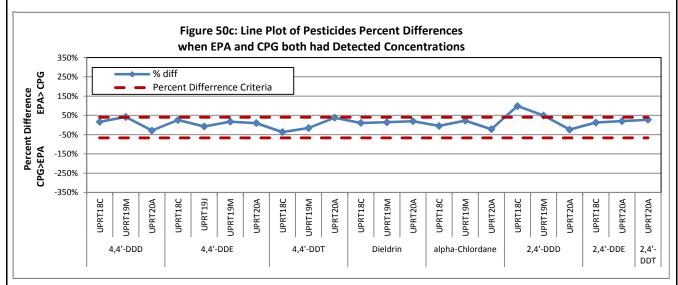
HpCDD = heptachlorodibenzo-p-dioxin TCDD = tetrachlorodibenzo-p-dioxin OCDD = octachlorodibenzo-p-dioxin

HpCDF = heptachlorodibenzofuran TCDF = tetrachlorodibenzofuran OCDF = octachlorodibenzofuran











Statistical Plot of Sediment Pesticides Scaled Concentrations

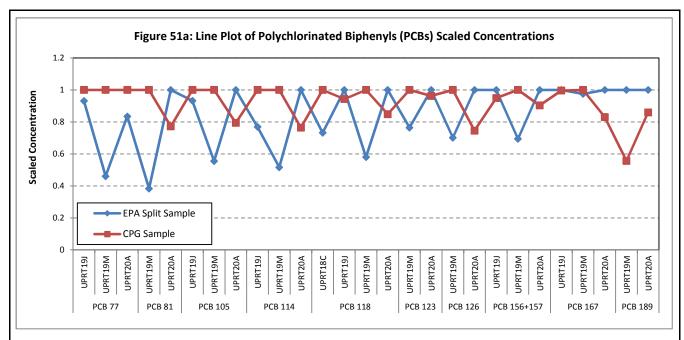
Figure 50

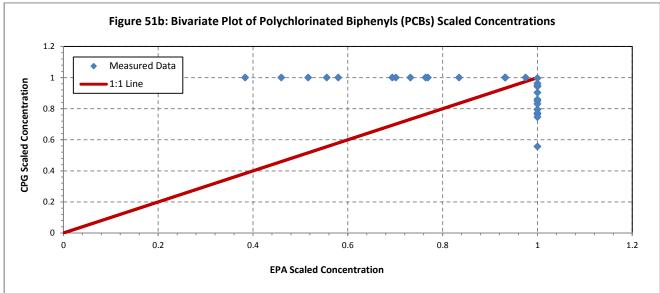
DDD = dichlorodiphenyldichloroethane

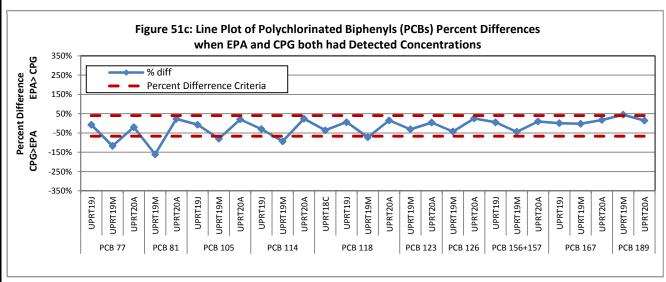
DDE = dichlorodiphenyldichloroethylene

DDT = dichlorodiphenyltrichloroethane







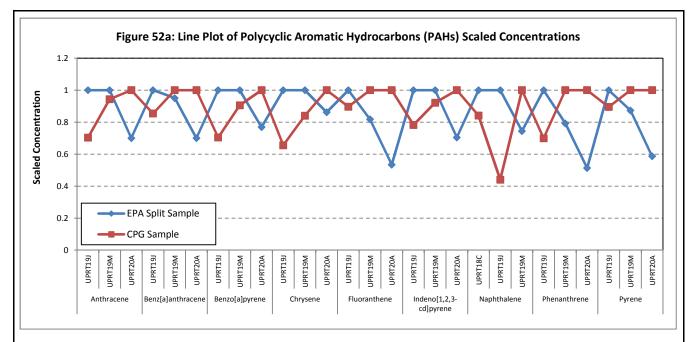


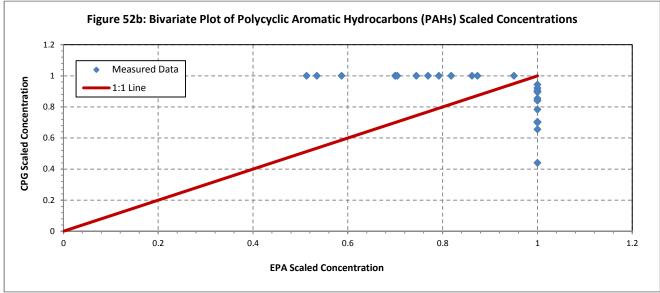


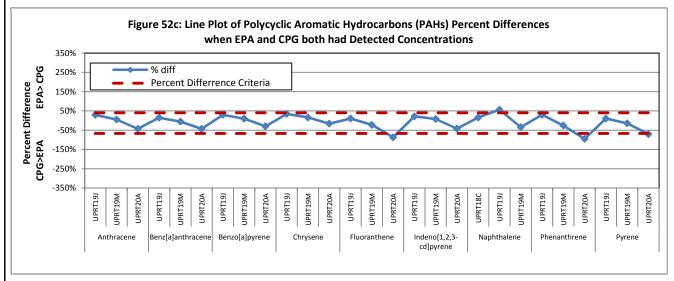
Statistical Plot of Sediment Polychlorinated Biphenyls (PCBs) Scaled Concentrations







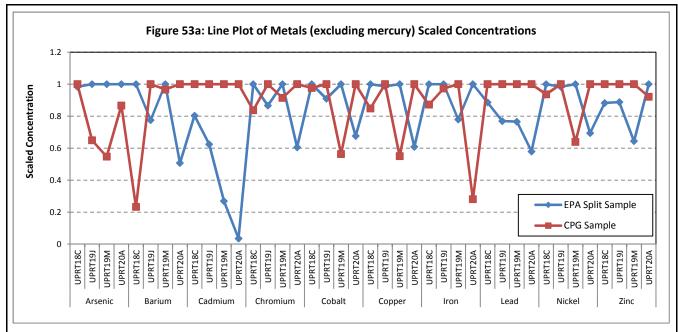


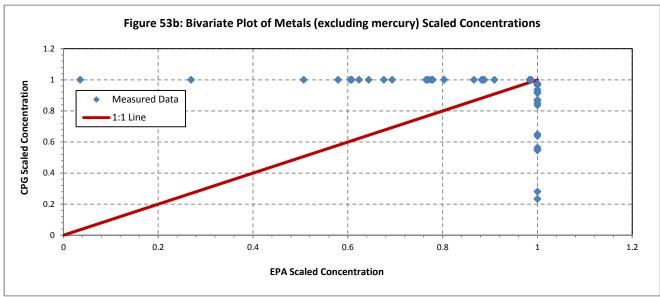


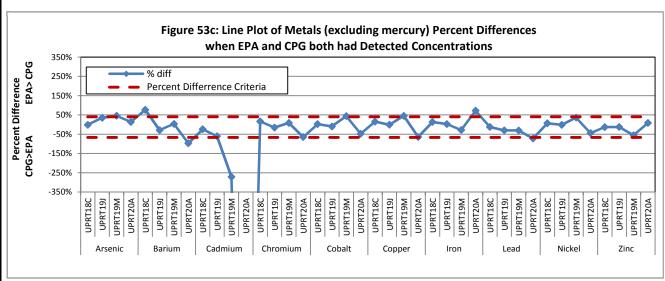


Statistical Plot of Sediment Polycyclic Aromatic Hydrocarbons (PAHs)
Scaled Concentrations











Statistical Plot of Sediment Metals (excluding mercury) Scaled Concentrations

